

Analytical Resources, Incorporated
Analytical Chemists and Consultants

October 14, 2011



John Long
AMEC/Geomatrix
600 University Suite 600
Seattle, WA 98101

RE: Client Project: Former Rhone Poulenc- 8769 Shoreline Investigation
ARI Job Numbers: TN42, TN43

Dear John:

Please find enclosed the final data package for samples for the project referenced above. ARI received six water samples and one trip blank on September 21, 2011.

Please refer to the case narrative for details on the analyses of these samples.

An electronic copy of this package will be kept on file at ARI. If you have questions or problems, please feel free to contact me at any time.

Sincerely,

ANALYTICAL RESOURCES, INC.

Cheronne Oreiro
Project Manager
-For-
Kelly Bottem
Client Services Manager
206/695-6211
kellyb@arilabs.com

Enclosures

cc: file TN42_TN43

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Project: 8769 FRP Shoreline Investigation

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(JL)
Signature

October-07-2011
Date

Chain of Custody Documentation

ARI Job ID: TN42, TN43

TN42 : 00002

Chain of Custody Record & Laboratory Analysis Request

Analytical Resources, Incorporated
Analytical Chemists and Consultants
4611 South 134th Place, Suite 100
Tukwila, WA 98168
206-695-6200 206-695-6201 (fax)

ARI Assigned Number:	1N47	Turn-around Requested:	Standard	Page:	1	of	1
ARI Client Company:	AMEC	Phone:	206-342-1779	Date:	9/21/11	Ce Present?	
Client Contact:	John Long / Nik Bacher	No. of Coolers:		No. of Coolers:	1	Cooler Temps:	2
Client Project Name:	FEP 2011 Shoreline Investigation	Analysis Requested					
Client Project #:	8769	Samplers:	Taylor Louvise	Date	Time	Matrix	No. Containers
Sample ID							
FEP-092111-001	9/21/11	0920	W	5			
FEP-092111-002		1110		5			
FEP-092111-003		1235		5			
FEP-092111-004		1300		5			
FEP-092111-005		1445		5			
FEP-092111-006		1600	V	5			
Trip Blank			W	2			
Comments/Special Instructions							
Metals include Al, As, Cd, Cr, Cu, Pb, Ni, Se, Tl, V, Zn.							
Received by: <i>The Trevor</i> (Signature)				Relinquished by: <i>Joe Wags</i> (Signature)			
Printed Name: Trevor Louvise Company: AMEC				Printed Name: Joe Wags Company: AMEC			
Date & Time:	9/21/11 1640	Date & Time:	9/21/11 16:40	Date & Time:	9/21/11	Date & Time:	9/21/11 16:40

Limits of Liability: API will perform all requested services in accordance with appropriate methodology following API Standard Operating Procedures and the API Quality Assurance Program. This program meets standards for the industry. The total liability of API, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by API release API from any liability in excess thereof, notwithstanding any provision to the contrary in any contract, purchase order or co-

Sample Retention Policy: All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.



Analytical Resources, Incorporated
Analytical Chemists and Consultants

Cooler Receipt Form

ARI Client: AMEC

COC No(s): _____ NA

Assigned ARI Job No: TN42

Preliminary Examination Phase:

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES NO

Were custody papers included with the cooler? YES NO

Were custody papers properly filled out (ink, signed, etc.) YES NO

Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry)..... 1.2

If cooler temperature is out of compliance fill out form 00070F

Temp Gun ID#: 90941617

Cooler Accepted by: JU (AN) Date: 9/21/11 Time: 1640

Complete custody forms and attach all shipping documents

Log-In Phase:

Was a temperature blank included in the cooler? YES NO

What kind of packing material was used? ... Bubble Wrap Wet Ice Gel Packs Baggies Foam Block Paper Other: Bubble Wrap Wet Ice Gel Packs Baggies Foam Block Paper Other: NA YES NO

Was sufficient ice used (if appropriate)? YES NO

Were all bottles sealed in individual plastic bags? YES NO

Did all bottles arrive in good condition (unbroken)? YES NO

Were all bottle labels complete and legible? YES NO

Did the number of containers listed on COC match with the number of containers received? YES NO

Did all bottle labels and tags agree with custody papers? YES NO

Were all bottles used correct for the requested analyses? YES NO

Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... NA YES NO

Were all VOC vials free of air bubbles? NA YES NO

Was sufficient amount of sample sent in each bottle? YES NO

Date VOC Trip Blank was made at ARI..... NA 9-19-11

Was Sample Split by ARI : NA YES Date/Time: _____ Equipment: _____ Split by: _____

Samples Logged by: TS Date: 9-21-11 Time: 1640

*** Notify Project Manager of discrepancies or concerns ***

Sample ID on Bottle	Sample ID on COC	Sample ID on Bottle	Sample ID on COC

Additional Notes, Discrepancies, & Resolutions:

FRP - 003 2"sm" FRP - 005 1"sm"
FRP - 004 1"sm"

By: TS

Date: 9-21-11

Small Air Bubbles ~2mm • * •	Peabubbles 2-4 mm • • •	LARGE Air Bubbles > 4 mm • • •	Small → "sm" Peabubbles → "pb" Large → "lg" Headspace → "hs"

PRESERVATION VERIFICATION 09/21/11

Page 1 of 1

Inquiry Number: NONE

Analysis Requested: 09/22/11

Contact: Long, John

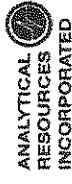
Client: AMEC Geomatix

Logged by: TS

Sample Set Used: Yes~481

Validatable Package: No

Deliverables:



ARI Job No: TN42

PC: Kelly

VTSR: 09/21/11

Project #: 8769
 Project: FRP 2011 Shoreline Investigation
 Sample Site:
 SDG No:
 Analytical Protocol: In-house

LOGNUM ARI ID	CLIENT ID	CN >12	WAD >12	NH3 <2	COD <2	FOG <2	MET <2	PHEN <2	TKN <2	NO23 <2	TOC <2	S2 <9	AR102 <2	Fe2+ <2	DMET FLT	DOC FLT	PARAMETER	ADJUSTED TO	LOT NUMBER	AMOUNT ADDED	DATE/BY
11-20663	FRP-092111-001																				
TN42A																					
11-20664	FRP-092111-002																				
TN42B																					
11-20665	FRP-092111-003																				
TN42C																					
11-20666	FRP-092111-004																				
TN42D																					
11-20667	FRP-092111-005																				
TN42E																					
11-20668	FRP-092111-006																				
TN42F																					

weak ph

X42 : 060205

TS
Checked By _____ Date _____



Analytical Resources, Incorporated
Analytical Chemists and Consultants

Cooler Receipt Form

ARI Client: Amec

Project Name: FRP 2011 Shoreline investigation

COC No(s): _____ NA

Delivered by: Fed-Ex UPS Courier Hand Delivered Other: _____

Assigned ARI Job No: TN43

Tracking No: _____ NA

Preliminary Examination Phase:

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES NO

Were custody papers included with the cooler? YES NO

Were custody papers properly filled out (ink, signed, etc.) YES NO

Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry) 12

If cooler temperature is out of compliance fill out form 00070F

Temp Gun ID#: 90941619

Cooler Accepted by: TS Date: 9-21-11 Time: 140

Complete custody forms and attach all shipping documents

Log-In Phase:

Was a temperature blank included in the cooler? YES NO

What kind of packing material was used? ... Bubble Wrap Wet Ice Gel Packs Baggies Foam Block Paper Other: _____ NA YES NO

Was sufficient ice used (if appropriate)? YES NO

Were all bottles sealed in individual plastic bags? YES NO

Did all bottles arrive in good condition (unbroken)? YES NO

Were all bottle labels complete and legible? YES NO

Did the number of containers listed on COC match with the number of containers received? YES NO

Did all bottle labels and tags agree with custody papers? YES NO

Were all bottles used correct for the requested analyses? YES NO

Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... NA YES NO

Were all VOC vials free of air bubbles? YES NO NA YES NO

Was sufficient amount of sample sent in each bottle? YES NO NA YES NO

Date VOC Trip Blank was made at ARI..... NA

Was Sample Split by ARI : NA YES Date/Time: _____ Equipment: _____ Split by: _____

Samples Logged by: TS Date: 9-21-11 Time: 1650

**** Notify Project Manager of discrepancies or concerns ****

Sample ID on Bottle	Sample ID on COC	Sample ID on Bottle	Sample ID on COC

Additional Notes, Discrepancies, & Resolutions:

By:

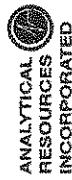
Date:

Small Air Bubbles ~2mm * * *	Peabubbles 2-4 mm * * *	LARGE Air Bubbles > 4 mm * * *	Small → "sm" Peabubbles → "pb" Large → "lg" Headspace → "hs"
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PRESERVATION VERIFICATION 09/21/11

Page 1 of 1

ARI Job No: TN43



Inquiry Number: NONE
 Analysis Requested: 09/22/11
 Contact: Long, John
 Client: AMEC Geomatrix
 Logged by: TS
 Sample Set Used: Yes-481
 Validatable Package: No
 Deliverables:

Project #: 8769
 Project: FRP Shoreline Investigation
 Sample Site:
 SDG No:
 Analytical Protocol: In-house

LOGNUM ARI ID	CLIENT ID	CN >12	WAD >12	NH3 <2	COD <2	FOG <2	MET <2	PHOS <2	TKN <2	NO23 <2	TOC <2	S2 >9	AK102 <2	Fe2+ <2	DMET FLT	DOC FLT	ADJUSTED PARAMETER	LCT TO	NUMBER	AMOUNT ADDED	DATE/BY
11-20670	FRP-092111-001																	1	11030	0.5mL	9.22.11 OM
11-20671	FRP-092111-002																	2	11030	0.5mL	9.22.11 OM
11-20672	FRP-092111-003																	2	11030	0.5mL	9.22.11 OM
11-20673	FRP-092111-004																	2	11030	0.5mL	9.22.11 OM
11-20674	FRP-092111-005																	1	11030	0.5mL	9.22.11 OM
11-20675	FRP-092111-006																	1	11030	0.5mL	9.22.11 OM

Wet lab pH

1842: 060007

TS
Checked By _____ Date _____

Case Narrative, Data Qualifiers, Control Limits

ARI Job ID: TN42, TN43

Case Narrative

AMEC/Geomatrix

Client Project: Former Rhone Poulenc- 8769 Shoreline Investigation

ARI Job Numbers: TN42, TN43

October 14, 2011

Page 1 of 2

Sample Receipt:

Please find enclosed the original Chain-of-Custody (COC) record, sample receipt documentation, and analytical results for the project referenced above. Analytical Resources, Inc. (ARI) accepted six water samples and one trip blank in good condition on 9/21/11. Please see the enclosed Cooler Receipt Form for further details.

Volatiles by 8260C

The samples were analyzed on 9/26/11 within the method recommended holding times.

Initial calibration (s): All analytes of interest were within method acceptance criteria.

Continuing calibration (s): The continuing calibration (CCAL) on 9/26/11 fell outside the 20% control limit low for Acetone, Methyl Iodide, and trans-1,4-Dichloro-2-butene. All detected results associated with this CCAL have been flagged with a "Q" qualifier.

LCS/LCSD/RPDs: The LCS percent recovery of 2-Chloroethylvinylether fell outside the control limits low for **LCS-092611**.

Surrogates: All surrogate recoveries were within control limits.

Method Blank (s): The method blank was free of contamination.

Samples: Samples **FRP-092111-001**, **FRP-092111-002**, **FRP-092111-003**, **FRP-092111-004**, **FRP-092111-005**, and **FRP-092111-006** had a pH between 6 and 8.

Metals Analysis (6010, 200.8 and 7000 series)

The samples were digested on 9/23/11 and 10/4/11 - within the method recommended holding time and analyzed between 9/27/11 and 10/7/11.

Initial calibration (s): All analytes of interest were within method acceptance criteria.

Continuing calibration (s): All analytes of interest were within method acceptance criteria.

LCS (s): The percent recoveries were within control limits.

Method Blank (s): The method blanks were free of contamination.

Samples: There were no anomalies associated with these samples.

Case Narrative

AMEC/Geomatrix

Client Project: Former Rhone Poulenc- 8769 Shoreline Investigation

ARI Job Numbers: TN42, TN43

October 14, 2011

Page 2 of 2

Matrix spike/ Sample Duplicate RPD(s): The matrix spike percent recovery of selenium fell outside the control limits high for sample FRP-092111-001. A post digestion spike was performed and the recovery was within control limits. All relevant data have been flagged with an "N" qualifier on the appropriate Form V.

The matrix spike percent recovery of mercury fell outside the control limits low for sample FRP-092111-001. All relevant data have been flagged with an "N" qualifier on the appropriate Form V.

The duplicate RPD of mercury was outside the control limit for sample FRP-092111-001. All relevant data have been flagged with a "*" qualifier on the appropriate Form VI.

pH by method 150.1:

The samples were analyzed on 9/21/11 within method recommended holding time.

LCS (s): The LCS percent recovery was within control limits.

Sample Replicate RPD (s): The RPD is in control.

Sample ID Cross Reference Report**ANALYTICAL
RESOURCES INCORPORATED**

ARI Job No: TN42

Client: AMEC Geomatrix

Project Event: 8769

Project Name: FRP 2011 Shoreline Investigation

Sample ID	ARI Lab ID	ARI LIMS ID	Matrix	Sample Date/Time	VTSR
1. FRP-092111-001	TN42A	11-20663	Water	09/21/11 09:20	09/21/11 16:40
2. FRP-092111-002	TN42B	11-20664	Water	09/21/11 11:10	09/21/11 16:40
3. FRP-092111-003	TN42C	11-20665	Water	09/21/11 12:55	09/21/11 16:40
4. FRP-092111-004	TN42D	11-20666	Water	09/21/11 13:00	09/21/11 16:40
5. FRP-092111-005	TN42E	11-20667	Water	09/21/11 14:45	09/21/11 16:40
6. FRP-092111-006	TN42F	11-20668	Water	09/21/11 16:00	09/21/11 16:40
7. Trip Blanks	TN42G	11-20669	Water	09/21/11	09/21/11 16:40

Printed 09/21/11

TN42 : 000012

Sample ID Cross Reference Report**ANALYTICAL
RESOURCES
INCORPORATED**

ARI Job No: TN43

Client: AMEC Geomatrix

Project Event: 8769

Project Name: FRP Shoreline Investigation

Sample ID	ARI Lab ID	ARI LIMS ID	Matrix	Sample Date/Time	VTSR
1. FRP-092111-001	TN43A	11-20670	Water	09/21/11 09:20	09/21/11 16:40
2. FRP-092111-002	TN43B	11-20671	Water	09/21/11 11:10	09/21/11 16:40
3. FRP-092111-003	TN43C	11-20672	Water	09/21/11 12:55	09/21/11 16:40
4. FRP-092111-004	TN43D	11-20673	Water	09/21/11 13:00	09/21/11 16:40
5. FRP-092111-005	TN43E	11-20674	Water	09/21/11 14:45	09/21/11 16:40
6. FRP-092111-006	TN43F	11-20675	Water	09/21/11 16:00	09/21/11 16:40

Printed 09/21/11



Analytical Resources, Incorporated
Analytical Chemists and Consultants

Data Reporting Qualifiers

Effective 2/14/2011

Inorganic Data

- U Indicates that the target analyte was not detected at the reported concentration
- * Duplicate RPD is not within established control limits
- B Reported value is less than the CRDL but \geq the Reporting Limit
- N Matrix Spike recovery not within established control limits
- NA Not Applicable, analyte not spiked
- H The natural concentration of the spiked element is so much greater than the concentration spiked that an accurate determination of spike recovery is not possible
- L Analyte concentration is \leq 5 times the Reporting Limit and the replicate control limit defaults to ± 1 RL instead of the normal 20% RPD

Organic Data

- U Indicates that the target analyte was not detected at the reported concentration
- * Flagged value is not within established control limits
- B Analyte detected in an associated Method Blank at a concentration greater than one-half of ARI's Reporting Limit or 5% of the regulatory limit or 5% of the analyte concentration in the sample.
- J Estimated concentration when the value is less than ARI's established reporting limits
- D The spiked compound was not detected due to sample extract dilution
- E Estimated concentration calculated for an analyte response above the valid instrument calibration range. A dilution is required to obtain an accurate quantification of the analyte.
- Q Indicates a detected analyte with an initial or continuing calibration that does not meet established acceptance criteria (<20%RSD, <20%Drift or minimum RRF).



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- S Indicates an analyte response that has saturated the detector. The calculated concentration is not valid; a dilution is required to obtain valid quantification of the analyte
- NA The flagged analyte was not analyzed for
- NR Spiked compound recovery is not reported due to chromatographic interference
- NS The flagged analyte was not spiked into the sample
- M Estimated value for an analyte detected and confirmed by an analyst but with low spectral match parameters. This flag is used only for GC-MS analyses
- M2 The sample contains PCB congeners that do not match any standard Aroclor pattern. The PCBs are identified and quantified as the Aroclor whose pattern most closely matches that of the sample. The reported value is an estimate.
- N The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification"
- Y The analyte is not detected at or above the reported concentration. The reporting limit is raised due to chromatographic interference. The Y flag is equivalent to the U flag with a raised reporting limit.
- EMPC Estimated Maximum Possible Concentration (EMPC) defined in EPA Statement of Work DLM02.2 as a value "calculated for 2,3,7,8-substituted isomers for which the quantitation and /or confirmation ion(s) has signal to noise in excess of 2.5, but does not meet identification criteria" **(Dioxin/Furan analysis only)**
- C The analyte was positively identified on only one of two chromatographic columns. Chromatographic interference prevented a positive identification on the second column
- P The analyte was detected on both chromatographic columns but the quantified values differ by ≥40% RPD with no obvious chromatographic interference
- X Analyte signal includes interference from polychlorinated diphenyl ethers. **(Dioxin/Furan analysis only)**
- Z Analyte signal includes interference from the sample matrix or perfluorokerosene ions. **(Dioxin/Furan analysis only)**



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Geotechnical Data

- A The total of all fines fractions. This flag is used to report total fines when **only** sieve analysis is requested and balances total grain size with sample weight.
- F Samples were frozen prior to particle size determination
- SM Sample matrix was not appropriate for the requested analysis. This normally refers to samples contaminated with an organic product that interferes with the sieving process and/or moisture content, porosity and saturation calculations
- SS Sample did not contain the proportion of "fines" required to perform the pipette portion of the grain size analysis
- W Weight of sample in some pipette aliquots was below the level required for accurate weighting

SURR SOLUTIONS

LABEL SOLN ID TEST CONC. UG/ML SOLVENT EXP.

LCS SOLUTIONS

LABL	SOLN ID	TEST	CONC.	UG/ML SOLVENT	EXP.
1	1888-2	PCB 1660	20	ACETONE	08/30/12
2#	NA	BCOC PEST	10	ACETONE	NA
3	1885-1	PEST	01/02/10	ACETONE	12/15/11
4	1885-2	LOW PEST	.1/.2/1	ACETONE	12/15/11
5	1779-1	EPH	1500	MECL2	11/11/11
6	1791-5	PCP	12.5/125	ACETONE	12/10/11
7	1888-1	ABN	100	MEOH	08/30/12
8	1785-3	TBT	2.5	MECL2	11/27/11
9	1786-3	PORE TBT	.125/.25	MECL2	11/27/11
10					
11	1860-4	TPHD	15000	ACETONE	05/12/12
12					
13	1838-4	LOW PCB	2	ACETONE	01/31/12
14					
15	1814-2	SIM PNA	15/75	MEOH	01/04/12
16	1879-3	1,4-DIOXANE	100	MEOH	02/05/12
17	1869-4	1248 PCB	10	ACETONE	06/14/12
18	1814-3	LOW SIM PNA	1.5	ACETONE	01/04/12
19	1873-2	AK103	7500	ACETONE	01/02/12
20	1886-4	PNA	100	ACETONE	01/07/12
21	1874-3	SKY/BHT	100	MEOH	01/14/12
22	1864-3	HERB	02 to 2500	MEOH	12/03/11
23	1887-2	EXTRA PNA	15	ACETONE	08/25/12
24					
25#	NA	DIPHENYL	100	MEOH	NA
26	1869-1	OP-PEST	25	MEOH	10/01/11
27	NA	STEROLS	200	MEOH	NA
28#	1807-1	ADD. PEST	2	ACETONE	08/31/11
29#	NA	DECANES	100	MEOH	NA

LCS SOLUTIONS

30	NA	EDB/DBCP	0.2	MEOH	NA
31	1835-2	TERPINEOL	100	MEOH	09/02/11
32	1876-1	GUAIACOL	50-200	ACETONE	01/05/12
33	NA	RETENE	100	MEOH	NA
34	1867-3	CONGENERS	0.5	ACETONE	03/14/12
35	1875-3	ALKYL PNA A	10	MEOH	07/18/12
36	NA	ALKYL PNA B	10	MEOH	NA
37	1773-1	CAR/PERY	100	ACETONE	10/14/11
38	1872-2	ABN ACID	200-450	MEOH	12/29/11
39	1853-4	BENZIDINE	500	MEOH	04/30/12
40	1851-3	PBDE	0.5	MEOH	04/22/12
50	1900-1	FULL RESIN	250	ACETONE	08/12/12
51	1772-1	DDTS	0.01	ACETONE	04/24/11
52	NA	1232 PCB	20	ACETONE	NA
53	1852-2	DALAPON	50	MEOH	12/03/11
54	1753-1	T-CHLORDANE	10	ACETONE	07/21/11
55	1753-2	TOXAPHENE	50	ACETONE	07/21/11
56	1874-1	ABN BASE	50-200	MEOH	01/05/12
#=PROJECT SPECIFIC SOLUTION					
*=REVERIFIED SOLUTION					



**Spike Recovery Control Limits for Analysis of Aqueous Samples
Volatile Organic Compounds (VOA) EPA SW-846 Methods 8260C
10 mL Purge Volume^(1,6)**

Effective: 8/30/2010

Control limits are updated periodically. Assure that you have ARI's current control limits by downloading the files at the time of use. http://www.arilabs.com/portal/downloads/ARI-CLs.zip		
	ARI Control Limits	ARI ME Control Limits ⁽²⁾
LCS Spike Recovery⁽⁵⁾		
Dichlorodifluoromethane	69 - 122	60 - 131
Chloromethane	76 - 120	69 - 123
Vinyl Chloride	80 - 120	75 - 123
Bromomethane	71 - 120	63 - 129
Chloroethane	80 - 120	75 - 124
1,1,2-Trichloro-1,2,2-trifluoroethane	80 - 121	76 - 128
Acrolein	69 - 126	60 - 136
Acetone	71 - 120	64 - 120
1,1-Dichloroethene	80 - 120	79 - 122
Bromoethane	80 - 120	80 - 121
Methyl Iodide	76 - 120	69 - 127
Methylene Chloride	80 - 120	77 - 120
Acrylonitrile	79 - 120	74 - 120
Methyl tert-Butyl Ether	80 - 120	77 - 121
Carbon Disulfide	80 - 120	78 - 121
trans-1,2-Dichloroethene	80 - 120	80 - 120
Vinyl Acetate	80 - 120	76 - 120
1,1-Dichloroethane	80 - 120	80 - 120
2-Butanone	80 - 120	76 - 120
2,2-Dichloropropane	80 - 120	77 - 120
cis-1,2-Dichloroethene	80 - 120	80 - 120
Chloroform	80 - 120	80 - 120
Bromodichloromethane	80 - 120	80 - 120
1,1,1-Trichloroethane	80 - 120	80 - 120
1,1-Dichloropropene	80 - 120	80 - 120
Carbon Tetrachloride	80 - 120	80 - 123
1,2-Dichloroethane	80 - 120	80 - 120
Benzene	80 - 120	80 - 120
Trichloroethene	80 - 120	80 - 120
1,2-Dichloropropane	80 - 120	80 - 120
Bromochloromethane	80 - 120	80 - 120
Dibromomethane	80 - 120	80 - 120
2-Chloroethylvinylether	80 - 120	75 - 120
4-Methyl-2-Pentanone	80 - 120	78 - 120
cis-1,3-Dichloropropene	80 - 120	80 - 120
Toluene	80 - 120	80 - 120
trans-1,3-Dichloropropene	80 - 120	80 - 120



2-Hexanone	80 - 120	75 - 120
1,1,2-Trichloroethane	80 - 120	80 - 120
1,3-Dichloropropane	80 - 120	80 - 120
Tetrachloroethene	80 - 120	80 - 120
Dibromochloromethane	80 - 120	80 - 120
Ethylene Dibromide	80 - 120	80 - 120
Chlorobenzene	80 - 120	80 - 120
Ethylbenzene	80 - 120	80 - 121
1,1,2,2-Tetrachloroethane	80 - 120	78 - 120
m,p-Xylene	80 - 120	80 - 120
o-Xylene	80 - 120	80 - 120
Styrene	80 - 120	80 - 122
Trichlorofluoromethane	80 - 120	78 - 123
Isopropylbenzene	80 - 120	79 - 121
Bromoform	80 - 120	79 - 120
1,1,1,2-Tetrachloroethane	80 - 120	80 - 120
1,2,3-Trichloropropane	80 - 120	77 - 120
trans-1,4-Dichloro-2-butene	74 - 122	66 - 130
n-Propylbenzene	80 - 120	80 - 120
Bromobenzene	80 - 120	78 - 120
1,3,5-Trimethylbenzene	80 - 120	80 - 120
2-Chlorotoluene	80 - 120	80 - 120
4-Chlorotoluene	80 - 120	80 - 120
tert-Butylbenzene	80 - 120	80 - 121
1,2,4-Trimethylbenzene	80 - 120	80 - 120
sec-Butylbenzene	80 - 120	80 - 121
4-Isopropyltoluene	80 - 120	80 - 123
1,3-Dichlorobenzene	80 - 120	80 - 120
1,4-Dichlorobenzene	80 - 120	80 - 120
n-Butylbenzene	80 - 120	80 - 122
1,2-Dichlorobenzene	80 - 120	80 - 120
1,2-Dibromo-3-chloropropane	76 - 120	71 - 120
1,2,4-Trichlorobenzene	77 - 120	71 - 120
Hexachloro-1,3-butadiene	77 - 120	70 - 127
Naphthalene	76 - 120	70 - 120
1,2,3-Trichlorobenzene	79 - 120	74 - 120
MB/LCS Surrogate Recovery		
Dibromofluoromethane	80 - 120	(3)
d4-1,2-Dichloroethane	80 - 120	(3)
d8-Toluene	80 - 120	(3)
4-Bromofluorobenzene	80 - 120	(3)
d4-1,2-Dichlorobenzene	80 - 120	(3)
 Sample Surrogate Recovery		
Dibromofluoromethane	80 - 120	(3)



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Analytical Chemists and Consultants

d4-1,2-Dichloroethane	80 - 120	(3)
d8-Toluene	80 - 120	(3)
4-Bromofluorobenzene	80 - 120	(3)
D4-1,2-Dichlorobenzene	80 - 120	(3)

(1) Control Limits calculated using all data generated 7/1/09 through 6/30/10.

(2) ME = A **marginal exceedance** defined in the NELAC Standard⁽⁴⁾ as beyond the LCS-CL but still within the ME limits. ME limits are between 3 and 4 standard deviations around the mean. A maximum of four marginal exceedances are acceptable. Five or more marginal exceedances require corrective action.

(3) Marginal Exceedances not allowed for surrogate standards. A corrective action is required for each surrogate recovery outside of the control limit range.

(4) **2003 NELAC Standard (EPA/600/R-04/003), July 2003**, Chapter 5, pages 251-252.

(5) Laboratory Control Sample (LCS) spike recovery control limits also used as advisory control limits for sample matrix spike (MS) analyzes. MS recovery values are advisory and not used to assess the acceptability of an analytical batch.

(6) Highlighted control limits (**bold font**) are adjusted from the calculated values as follows:

a) ARI does not use control limits < 10 for the lower limit or < 100 for the upper limit.

b) Control limits for analyzes with no separate preparation procedure are adjusted to reflect the minimum uncertainty in the calibration of the instrument allowed by the referenced analytical method.



Summary of Laboratory Control Limits Metals Analyses (All Methods & Sample Matrices)

Effective 5/1/09

Control limits are updated periodically. Assure that you have ARI's current control limits by downloading the files at the time of use. <http://www.arilabs.com/portal/downloads/ARI-CLs.zip>

Element	Matrix Spike Recovery	LCS Recovery	Replicate RPD
Aluminum	75 - 125	80 - 120	≤ 20%
Antimony	75 - 125	80 - 120	≤ 20%
Arsenic	75 - 125	80 - 120	≤ 20%
Barium	75 - 125	80 - 120	≤ 20%
Beryllium	75 - 125	80 - 120	≤ 20%
Boron	75 - 125	80 - 120	≤ 20%
Cadmium	75 - 125	80 - 120	≤ 20%
Calcium	75 - 125	80 - 120	≤ 20%
Chromium	75 - 125	80 - 120	≤ 20%
Cobalt	75 - 125	80 - 120	≤ 20%
Copper	75 - 125	80 - 120	≤ 20%
Iron	75 - 125	80 - 120	≤ 20%
Lead	75 - 125	80 - 120	≤ 20%
Magnesium	75 - 125	80 - 120	≤ 20%
Manganese	75 - 125	80 - 120	≤ 20%
Mercury	75 - 125	80 - 120	≤ 20%
Nickel	75 - 125	80 - 120	≤ 20%
Potassium	75 - 125	80 - 120	≤ 20%
Selenium	75 - 125	80 - 120	≤ 20%
Silica	75 - 125	80 - 120	≤ 20%
Silver	75 - 125	80 - 120	≤ 20%
Sodium	75 - 125	80 - 120	≤ 20%
Strontium	75 - 125	80 - 120	≤ 20%
Thallium	75 - 125	80 - 120	≤ 20%
Vanadium	75 - 125	80 - 120	≤ 20%
Zinc	75 - 125	80 - 120	≤ 20%



Spike Recovery Control Limits for Conventional Wet Chemistry

Effective 5/1/09

Control limits are updated periodically. Assure that you have ARI's current control limits by downloading the files at the time of use. <http://www.arilabs.com/portal/downloads/ARI-CLs.zip>

Sample Matrix:	ARI's Control Limits	
	Water	Soil / Sediment
Matrix Spike Recoveries	% Recovery	% Recovery
Ammonia	75 - 125	75 - 125
Bromide	75 - 125	75 - 125
Chloride	75 - 125	75 - 125
Cyanide	75 - 125	75 - 125
Ferrous Iron	75 - 125	75 - 125
Fluoride	75 - 125	75 - 125
Formaldehyde	75 - 125	75 - 125
Hexane Extractable Material	-- -- --	78 - 114
Hexavalent Chromium	75 - 125	75 - 125
Nitrate/Nitrite	75 - 125	75 - 125
Oil and Grease	75 - 125	75 - 125
Phenol	75 - 125	75 - 125
Phosphorous	75 - 125	75 - 125
Sulfate	75 - 125	75 - 125
Sulfide	75 - 125	75 - 125
Total Kjeldahl Nitrogen	75 - 125	75 - 125
Total Organic Carbon	75 - 125	75 - 125
Duplicate RPDs		
Acidity	±20%	±20%
Alkalinity	±20%	±20%
BOD	±20%	±20%
Cation Exchange	±20%	±20%
COD	±20%	±20%
Conductivity	±20%	±20%
Salinity	±20%	±20%
Solids	±20%	±20%
Turbidity	±20%	±20%

**Volatile Analysis
Report and Summary QC Forms**

ARI Job ID: TN42, TN43

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2

Sample ID: FRP-092111-001
SAMPLE

Lab Sample ID: TN42A
LIMS ID: 11-20663
Matrix: Water
Data Release Authorized: *TMW*
Reported: 10/03/11

QC Report No: TN42-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769
Date Sampled: 09/21/11
Date Received: 09/21/11

Instrument/Analyst: NT2/PAB
Date Analyzed: 09/26/11 16:03

Sample Amount: 10.0 mL
Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U
75-15-0	Carbon Disulfide	0.09	0.2	0.3
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	0.3
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	3.5
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoro	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	<i>o</i> -Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

Sample ID: FRP-092111-001
SAMPLE

Lab Sample ID: TN42A
LIMS ID: 11-20663
Matrix: Water
Date Analyzed: 09/26/11 16:03

QC Report No: TN42-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromo-chloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	95.2%
d8-Toluene	99.5%
Bromo-fluorobenzene	94.4%
d4-1,2-Dichlorobenzene	101%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C

Page 1 of 2

**ANALYTICAL
RESOURCES
INCORPORATED**


Sample ID: FRP-092111-002

SAMPLE

Lab Sample ID: TN42B

LIMS ID: 11-20664

Matrix: Water

Data Release Authorized: *MW*

Reported: 10/03/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/21/11

Date Received: 09/21/11

Instrument/Analyst: NT2/PAB

Date Analyzed: 09/26/11 16:30

Sample Amount: 2.00 mL

Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.49	2.5	< 2.5 U
74-83-9	Bromomethane	0.22	5.0	< 5.0 U
75-01-4	Vinyl Chloride	0.38	1.0	< 1.0 U
75-00-3	Chloroethane	0.76	1.0	< 1.0 U
75-09-2	Methylene Chloride	2.0	2.5	< 2.5 U
67-64-1	Acetone	3.6	25	< 25 U
75-15-0	Carbon Disulfide	0.44	1.0	< 1.0 U
75-35-4	1,1-Dichloroethene	0.46	1.0	< 1.0 U
75-34-3	1,1-Dichloroethane	0.26	1.0	< 1.0 U
156-60-5	trans-1,2-Dichloroethene	0.42	1.0	< 1.0 U
156-59-2	cis-1,2-Dichloroethene	0.50	1.0	< 1.0 U
67-66-3	Chloroform	0.40	1.0	< 1.0 U
107-06-2	1,2-Dichloroethane	0.38	1.0	< 1.0 U
78-93-3	2-Butanone	4.0	25	< 25 U
71-55-6	1,1,1-Trichloroethane	0.44	1.0	< 1.0 U
56-23-5	Carbon Tetrachloride	0.38	1.0	< 1.0 U
108-05-4	Vinyl Acetate	0.34	5.0	< 5.0 U
75-27-4	Bromodichloromethane	0.26	1.0	< 1.0 U
78-87-5	1,2-Dichloropropane	0.46	1.0	< 1.0 U
10061-01-5	cis-1,3-Dichloropropene	0.29	1.0	< 1.0 U
79-01-6	Trichloroethene	0.38	1.0	< 1.0 U
124-48-1	Dibromochloromethane	0.45	1.0	< 1.0 U
79-00-5	1,1,2-Trichloroethane	0.18	1.0	< 1.0 U
71-43-2	Benzene	0.28	1.0	< 1.0 U
10061-02-6	trans-1,3-Dichloropropene	0.30	1.0	< 1.0 U
110-75-8	2-Chloroethylvinylether	0.43	5.0	< 5.0 U
75-25-2	Bromoform	0.35	1.0	< 1.0 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	1.9	25	< 25 U
591-78-6	2-Hexanone	1.6	25	< 25 U
127-18-4	Tetrachloroethene	0.44	1.0	< 1.0 U
79-34-5	1,1,2-Tetrachloroethane	0.34	1.0	< 1.0 U
108-88-3	Toluene	0.28	1.0	< 1.0 U
108-90-7	Chlorobenzene	0.21	1.0	< 1.0 U
100-41-4	Ethylbenzene	0.47	1.0	< 1.0 U
100-42-5	Styrene	0.33	1.0	< 1.0 U
75-69-4	Trichlorofluoromethane	0.46	1.0	< 1.0 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.54	1.0	< 1.0 U
179601-23-1	m,p-Xylene	0.72	2.0	< 2.0 U
95-47-6	o-Xylene	0.28	1.0	< 1.0 U
95-50-1	1,2-Dichlorobenzene	0.28	1.0	< 1.0 U
541-73-1	1,3-Dichlorobenzene	0.20	1.0	< 1.0 U
106-46-7	1,4-Dichlorobenzene	0.28	1.0	< 1.0 U
107-02-8	Acrolein	1.5	25	< 25 U
74-88-4	Methyl Iodide	0.20	5.0	< 5.0 U
74-96-4	Bromoethane	0.45	1.0	< 1.0 U
107-13-1	Acrylonitrile	0.92	5.0	< 5.0 U
563-58-6	1,1-Dichloropropene	0.46	1.0	< 1.0 U
74-95-3	Dibromomethane	0.40	1.0	< 1.0 U
630-20-6	1,1,1,2-Tetrachloroethane	0.34	1.0	< 1.0 U
96-12-8	1,2-Dibromo-3-chloropropane	1.1	2.5	< 2.5 U

MAP

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
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Sample ID: FRP-092111-002
SAMPLE

Lab Sample ID: TN42B
LIMS ID: 11-20664
Matrix: Water
Date Analyzed: 09/26/11 16:30

QC Report No: TN42-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	1.1	2.5	< 2.5 U
110-57-6	trans-1,4-Dichloro-2-butene	1.2	5.0	< 5.0 U
108-67-8	1,3,5-Trimethylbenzene	0.32	1.0	< 1.0 U
95-63-6	1,2,4-Trimethylbenzene	0.29	1.0	< 1.0 U
87-68-3	Hexachlorobutadiene	0.56	2.5	< 2.5 U
106-93-4	Ethylene Dibromide	0.38	1.0	< 1.0 U
74-97-5	Bromo-chloromethane	0.34	1.0	< 1.0 U
594-20-7	2,2-Dichloropropane	0.42	1.0	< 1.0 U
142-28-9	1,3-Dichloropropane	0.10	1.0	< 1.0 U
98-82-8	Isopropylbenzene	0.31	1.0	< 1.0 U
103-65-1	n-Propylbenzene	0.40	1.0	< 1.0 U
108-86-1	Bromobenzene	0.26	1.0	< 1.0 U
95-49-8	2-Chlorotoluene	0.21	1.0	< 1.0 U
106-43-4	4-Chlorotoluene	0.36	1.0	< 1.0 U
98-06-6	tert-Butylbenzene	0.30	1.0	< 1.0 U
135-98-8	sec-Butylbenzene	0.38	1.0	< 1.0 U
99-87-6	4-Isopropyltoluene	0.38	1.0	< 1.0 U
104-51-8	n-Butylbenzene	0.54	1.0	< 1.0 U
120-82-1	1,2,4-Trichlorobenzene	0.50	2.5	< 2.5 U
91-20-3	Naphthalene	0.35	2.5	< 2.5 U
87-61-6	1,2,3-Trichlorobenzene	0.44	2.5	< 2.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	93.8%
d8-Toluene	99.4%
Bromo-fluorobenzene	94.3%
d4-1,2-Dichlorobenzene	101%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2
**ANALYTICAL
RESOURCES
INCORPORATED**

Sample ID: FRP-092111-003
SAMPLELab Sample ID: TN42C
LIMS ID: 11-20665
Matrix: Water
Data Release Authorized: MW
Reported: 10/03/11QC Report No: TN42-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769
Date Sampled: 09/21/11
Date Received: 09/21/11Instrument/Analyst: NT2/PAB
Date Analyzed: 09/26/11 16:57Sample Amount: 10.0 mL
Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

MW

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
 Page 2 of 2

Sample ID: FRP-092111-003
 SAMPLE

Lab Sample ID: TN42C
 LIMS ID: 11-20665
 Matrix: Water
 Date Analyzed: 09/26/11 16:57

QC Report No: TN42-AMEC Geomatrix
 Project: FRP 2011 Shoreline Investigation
 8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromo-chloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	95.1%
d8-Toluene	99.3%
Bromo-fluorobenzene	94.3%
d4-1,2-Dichlorobenzene	101%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2

Sample ID: FRP-092111-004

SAMPLE

Lab Sample ID: TN42D
 LIMS ID: 11-20666
 Matrix: Water
 Data Release Authorized: MW
 Reported: 10/03/11

QC Report No: TN42-AMEC Geomatrix
 Project: FRP 2011 Shoreline Investigation
 8769
 Date Sampled: 09/21/11
 Date Received: 09/21/11

Instrument/Analyst: NT2/PAB
 Date Analyzed: 09/26/11 17:23

Sample Amount: 10.0 mL
 Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

Sample ID: FRP-092111-004
SAMPLE

Lab Sample ID: TN42D
LIMS ID: 11-20666
Matrix: Water
Date Analyzed: 09/26/11 17:23

QC Report No: TN42-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1, 2, 3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1, 4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1, 3, 5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1, 2, 4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromo-chloromethane	0.07	0.2	< 0.2 U
594-20-7	2, 2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1, 3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1, 2, 4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1, 2, 3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1, 2-Dichloroethane	94.1%
d8-Toluene	98.8%
Bromo-fluorobenzene	93.5%
d4-1, 2-Dichlorobenzene	102%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

ORGANICS ANALYSIS DATA SHEET

Volatile s by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2ANALYTICAL
RESOURCES
INCORPORATEDSample ID: FRP-092111-005
SAMPLE

Lab Sample ID: TN42E

LIMS ID: 11-20667

Matrix: Water

Data Release Authorized: *MMW*

Reported: 10/03/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/21/11

Date Received: 09/21/11

Instrument/Analyst: NT2/PAB

Date Analyzed: 09/26/11 17:50

Sample Amount: 2.00 mL

Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.49	2.5	< 2.5 U
74-83-9	Bromomethane	0.22	5.0	< 5.0 U
75-01-4	Vinyl Chloride	0.38	1.0	< 1.0 U
75-00-3	Chloroethane	0.76	1.0	< 1.0 U
75-09-2	Methylene Chloride	2.0	2.5	< 2.5 U
67-64-1	Acetone	3.6	25	< 25 U
75-15-0	Carbon Disulfide	0.44	1.0	< 1.0 U
75-35-4	1,1-Dichloroethene	0.46	1.0	< 1.0 U
75-34-3	1,1-Dichloroethane	0.26	1.0	< 1.0 U
156-60-5	trans-1,2-Dichloroethene	0.42	1.0	< 1.0 U
156-59-2	cis-1,2-Dichloroethene	0.50	1.0	< 1.0 U
67-66-3	Chloroform	0.40	1.0	< 1.0 U
107-06-2	1,2-Dichloroethane	0.38	1.0	< 1.0 U
78-93-3	2-Butanone	4.0	25	< 25 U
71-55-6	1,1,1-Trichloroethane	0.44	1.0	< 1.0 U
56-23-5	Carbon Tetrachloride	0.38	1.0	< 1.0 U
108-05-4	Vinyl Acetate	0.34	5.0	< 5.0 U
75-27-4	Bromodichloromethane	0.26	1.0	< 1.0 U
78-87-5	1,2-Dichloropropane	0.46	1.0	< 1.0 U
10061-01-5	cis-1,3-Dichloropropene	0.29	1.0	< 1.0 U
79-01-6	Trichloroethene	0.38	1.0	< 1.0 U
124-48-1	Dibromochloromethane	0.45	1.0	< 1.0 U
79-00-5	1,1,2-Trichloroethane	0.18	1.0	< 1.0 U
71-43-2	Benzene	0.28	1.0	< 1.0 U
10061-02-6	trans-1,3-Dichloropropene	0.30	1.0	< 1.0 U
110-75-8	2-Chloroethylvinylether	0.43	5.0	< 5.0 U
75-25-2	Bromoform	0.35	1.0	< 1.0 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	1.9	25	< 25 U
591-78-6	2-Hexanone	1.6	25	< 25 U
127-18-4	Tetrachloroethene	0.44	1.0	< 1.0 U
79-34-5	1,1,2,2-Tetrachloroethane	0.34	1.0	< 1.0 U
108-88-3	Toluene	0.28	1.0	< 1.0 U
108-90-7	Chlorobenzene	0.21	1.0	< 1.0 U
100-41-4	Ethylbenzene	0.47	1.0	< 1.0 U
100-42-5	Styrene	0.33	1.0	< 1.0 U
75-69-4	Trichlorofluoromethane	0.46	1.0	< 1.0 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoro	0.54	1.0	< 1.0 U
179601-23-1	m,p-Xylene	0.72	2.0	< 2.0 U
95-47-6	o-Xylene	0.28	1.0	< 1.0 U
95-50-1	1,2-Dichlorobenzene	0.28	1.0	< 1.0 U
541-73-1	1,3-Dichlorobenzene	0.20	1.0	< 1.0 U
106-46-7	1,4-Dichlorobenzene	0.28	1.0	< 1.0 U
107-02-8	Acrolein	1.5	25	< 25 U
74-88-4	Methyl Iodide	0.20	5.0	< 5.0 U
74-96-4	Bromoethane	0.45	1.0	< 1.0 U
107-13-1	Acrylonitrile	0.92	5.0	< 5.0 U
563-58-6	1,1-Dichloropropene	0.46	1.0	< 1.0 U
74-95-3	Dibromomethane	0.40	1.0	< 1.0 U
630-20-6	1,1,1,2-Tetrachloroethane	0.34	1.0	< 1.0 U
96-12-8	1,2-Dibromo-3-chloropropane	1.1	2.5	< 2.5 U

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2ANALYTICAL
RESOURCES
INCORPORATEDSample ID: FRP-092111-005
SAMPLELab Sample ID: TN42E
LIMS ID: 11-20667
Matrix: Water
Date Analyzed: 09/26/11 17:50QC Report No: TN42-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	1.1	2.5	< 2.5 U
110-57-6	trans-1,4-Dichloro-2-butene	1.2	5.0	< 5.0 U
108-67-8	1,3,5-Trimethylbenzene	0.32	1.0	< 1.0 U
95-63-6	1,2,4-Trimethylbenzene	0.29	1.0	< 1.0 U
87-68-3	Hexachlorobutadiene	0.56	2.5	< 2.5 U
106-93-4	Ethylene Dibromide	0.38	1.0	< 1.0 U
74-97-5	Bromoform	0.34	1.0	< 1.0 U
594-20-7	2,2-Dichloropropane	0.42	1.0	< 1.0 U
142-28-9	1,3-Dichloropropane	0.10	1.0	< 1.0 U
98-82-8	Isopropylbenzene	0.31	1.0	< 1.0 U
103-65-1	n-Propylbenzene	0.40	1.0	< 1.0 U
108-86-1	Bromobenzene	0.26	1.0	< 1.0 U
95-49-8	2-Chlorotoluene	0.21	1.0	< 1.0 U
106-43-4	4-Chlorotoluene	0.36	1.0	< 1.0 U
98-06-6	tert-Butylbenzene	0.30	1.0	< 1.0 U
135-98-8	sec-Butylbenzene	0.38	1.0	< 1.0 U
99-87-6	4-Isopropyltoluene	0.38	1.0	< 1.0 U
104-51-8	n-Butylbenzene	0.54	1.0	< 1.0 U
120-82-1	1,2,4-Trichlorobenzene	0.50	2.5	< 2.5 U
91-20-3	Naphthalene	0.35	2.5	< 2.5 U
87-61-6	1,2,3-Trichlorobenzene	0.44	2.5	< 2.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	93.2%
d8-Toluene	98.5%
Bromofluorobenzene	93.7%
d4-1,2-Dichlorobenzene	101%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.



ORGANICS ANALYSIS DATA SHEET

Volatile s by Purge & Trap GC/MS-Method SW8260C

Page 1 of 2

**ANALYTICAL
RESOURCES
INCORPORATED**


Sample ID: FRP-092111-006

SAMPLE

Lab Sample ID: TN42F

LIMS ID: 11-20668

Matrix: Water

Data Release Authorized: *MM*

Reported: 10/03/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/21/11

Date Received: 09/21/11

Instrument/Analyst: NT2/PAB

Date Analyzed: 09/26/11 18:17

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	7.4 <i>UJ</i>
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	1.1
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 <i>UJ</i>
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2
**ANALYTICAL
RESOURCES
INCORPORATED**


Sample ID: FRP-092111-006

SAMPLE

Lab Sample ID: TN42F
 LIMS ID: 11-20668
 Matrix: Water
 Date Analyzed: 09/26/11 18:17

QC Report No: TN42-AMEC Geomatrix
 Project: FRP 2011 Shoreline Investigation
 8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U J
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromo-chloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	92.3%
d8-Toluene	100%
Bromo-fluorobenzene	95.9%
d4-1,2-Dichlorobenzene	100%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

W/MH

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2Sample ID: Trip Blanks
SAMPLELab Sample ID: TN42G
LIMS ID: 11-20669
Matrix: Water
Data Release Authorized: *MW*
Reported: 10/03/11QC Report No: TN42-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769Date Sampled: 09/21/11
Date Received: 09/21/11Instrument/Analyst: NT2/PAB
Date Analyzed: 09/26/11 12:28Sample Amount: 10.0 mL
Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoro	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

Sample ID: Trip Blanks
SAMPLE

Lab Sample ID: TN42G
LIMS ID: 11-20669
Matrix: Water
Date Analyzed: 09/26/11 12:28

QC Report No: TN42-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromochloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in µg/L (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	94.1%
d8-Toluene	99.5%
Bromofluorobenzene	96.4%
d4-1,2-Dichlorobenzene	101%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

VOA SURROGATE RECOVERY SUMMARY

**ANALYTICAL
RESOURCES
INCORPORATED**



Matrix: Water

QC Report No: TN42-AMEC Geomatrix
 Project: FRP 2011 Shoreline Investigation
 8769

<u>ARI ID</u>	<u>Client ID</u>	<u>PV</u>	<u>DCE</u>	<u>TOL</u>	<u>BFB</u>	<u>DCB</u>	<u>TOT OUT</u>
MB-092611	Method Blank	10	94.0%	99.0%	96.0%	100%	0
LCS-092611	Lab Control	10	93.9%	98.3%	98.1%	101%	0
LCSD-092611	Lab Control Dup	10	93.3%	99.9%	98.3%	99.9%	0
TN42A	FRP-092111-001	10	95.2%	99.5%	94.4%	101%	0
TN42B	FRP-092111-002	10	93.8%	99.4%	94.3%	101%	0
TN42C	FRP-092111-003	10	95.1%	99.3%	94.3%	101%	0
TN42D	FRP-092111-004	10	94.1%	98.8%	93.5%	102%	0
TN42E	FRP-092111-005	10	93.2%	98.5%	93.7%	101%	0
TN42F	FRP-092111-006	10	92.3%	100%	95.9%	100%	0
TN42G	Trip Blanks	10	94.1%	99.5%	96.4%	101%	0

LCS/MB LIMITS QC LIMITS**SW8260C**

(DCE) = d4-1,2-Dichloroethane	80-120	80-120
(TOL) = d8-Toluene	80-120	80-120
(BFB) = Bromofluorobenzene	80-120	80-120
(DCB) = d4-1,2-Dichlorobenzene	80-120	80-120

Prep Method: SW5030B
 Log Number Range: 11-20663 to 11-20669

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2ANALYTICAL
RESOURCES
INCORPORATED

Sample ID: LCS-092611

LAB CONTROL SAMPLE

Lab Sample ID: LCS-092611

LIMS ID: 11-20663

Matrix: Water

Data Release Authorized: *MW*

Reported: 10/03/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: NA

Date Received: NA

Instrument/Analyst LCS: NT2/PAB

Sample Amount LCS: 10.0 mL

LCSD: NT2/PAB

LCSD: 10.0 mL

Date Analyzed LCS: 09/26/11 10:06

Purge Volume LCS: 10.0 mL

LCSD: 09/26/11 10:33

LCSD: 10.0 mL

Analyte	LCS	Spike Added-LCS	LCS Recovery	LCSD	Spike Added-LCSD	LCSD Recovery	RPD
Chloromethane	8.4	10.0	84.0%	8.9	10.0	89.0%	5.8%
Bromomethane	8.0	10.0	80.0%	8.3	10.0	83.0%	3.7%
Vinyl Chloride	9.7	10.0	97.0%	9.7	10.0	97.0%	0.0%
Chloroethane	8.8	10.0	88.0%	9.1	10.0	91.0%	3.4%
Methylene Chloride	9.2	10.0	92.0%	9.5	10.0	95.0%	3.2%
Acetone	41.3 Q	50.0	82.6%	41.9 Q	50.0	83.8%	1.4%
Carbon Disulfide	9.5	10.0	95.0%	9.6	10.0	96.0%	1.0%
1,1-Dichloroethene	8.8	10.0	88.0%	9.2	10.0	92.0%	4.4%
1,1-Dichloroethane	9.2	10.0	92.0%	9.3	10.0	93.0%	1.1%
trans-1,2-Dichloroethene	9.5	10.0	95.0%	9.5	10.0	95.0%	0.0%
cis-1,2-Dichloroethene	9.4	10.0	94.0%	9.7	10.0	97.0%	3.1%
Chloroform	9.5	10.0	95.0%	9.7	10.0	97.0%	2.1%
1,2-Dichloroethane	9.1	10.0	91.0%	9.4	10.0	94.0%	3.2%
2-Butanone	50.0	50.0	100%	49.8	50.0	99.6%	0.4%
1,1,1-Trichloroethane	9.4	10.0	94.0%	9.6	10.0	96.0%	2.1%
Carbon Tetrachloride	9.7	10.0	97.0%	9.9	10.0	99.0%	2.0%
Vinyl Acetate	9.6	10.0	96.0%	9.8	10.0	98.0%	2.1%
Bromodichloromethane	9.6	10.0	96.0%	9.8	10.0	98.0%	2.1%
1,2-Dichloropropane	9.3	10.0	93.0%	9.5	10.0	95.0%	2.1%
cis-1,3-Dichloropropene	9.6	10.0	96.0%	9.8	10.0	98.0%	2.1%
Trichloroethene	9.7	10.0	97.0%	9.9	10.0	99.0%	2.0%
Dibromochloromethane	10.0	10.0	100%	10.0	10.0	100%	0.0%
1,1,2-Trichloroethane	9.6	10.0	96.0%	9.9	10.0	99.0%	3.1%
Benzene	9.4	10.0	94.0%	9.8	10.0	98.0%	4.2%
trans-1,3-Dichloropropene	9.6	10.0	96.0%	9.6	10.0	96.0%	0.0%
2-Chloroethylvinylether	7.7	10.0	77.0%	8.3	10.0	83.0%	7.5%
Bromoform	10.4	10.0	104%	10.3	10.0	103%	1.0%
4-Methyl-2-Pentanone (MIBK)	49.8	50.0	99.6%	51.4	50.0	103%	3.2%
2-Hexanone	49.1	50.0	98.2%	49.6	50.0	99.2%	1.0%
Tetrachloroethene	9.9	10.0	99.0%	9.9	10.0	99.0%	0.0%
1,1,2,2-Tetrachloroethane	10.0	10.0	100%	9.9	10.0	99.0%	1.0%
Toluene	9.6	10.0	96.0%	9.9	10.0	99.0%	3.1%
Chlorobenzene	9.5	10.0	95.0%	9.7	10.0	97.0%	2.1%
Ethylbenzene	9.5	10.0	95.0%	9.5	10.0	95.0%	0.0%
Styrene	10.0	10.0	100%	10.2	10.0	102%	2.0%
Trichlorofluoromethane	9.7	10.0	97.0%	9.8	10.0	98.0%	1.0%
1,1,2-Trichloro-1,2,2-trifluoroethane	9.6	10.0	96.0%	9.7	10.0	97.0%	1.0%
m,p-Xylene	19.5	20.0	97.5%	19.8	20.0	99.0%	1.5%
o-Xylene	9.6	10.0	96.0%	9.7	10.0	97.0%	1.0%
1,2-Dichlorobenzene	9.6	10.0	96.0%	9.7	10.0	97.0%	1.0%
1,3-Dichlorobenzene	9.7	10.0	97.0%	9.7	10.0	97.0%	0.0%
1,4-Dichlorobenzene	9.6	10.0	96.0%	9.7	10.0	97.0%	1.0%
Acrolein	53.6	50.0	107%	56.1	50.0	112%	4.6%
Methyl Iodide	7.6 Q	10.0	76.0%	8.0 Q	10.0	80.0%	5.1%
Bromoethane	9.5	10.0	95.0%	9.6	10.0	96.0%	1.0%

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2
**ANALYTICAL
RESOURCES
INCORPORATED**

Sample ID: LCS-092611
LAB CONTROL SAMPLELab Sample ID: LCS-092611
LIMS ID: 11-20663
Matrix: WaterQC Report No: TN42-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

Analyte	LCS	Spike	LCS	Spike	LCSD	RPD
		Added-LCS	Recovery	LCSD	Added-LCSD	
Acrylonitrile	9.4	10.0	94.0%	9.7	10.0	97.0% 3.1%
1,1-Dichloropropene	9.6	10.0	96.0%	9.9	10.0	99.0% 3.1%
Dibromomethane	9.4	10.0	94.0%	9.7	10.0	97.0% 3.1%
1,1,1,2-Tetrachloroethane	9.9	10.0	99.0%	9.9	10.0	99.0% 0.0%
1,2-Dibromo-3-chloropropane	9.6	10.0	96.0%	9.4	10.0	94.0% 2.1%
1,2,3-Trichloropropane	9.5	10.0	95.0%	9.6	10.0	96.0% 1.0%
trans-1,4-Dichloro-2-butene	7.8 Q	10.0	78.0%	7.7 Q	10.0	77.0% 1.3%
1,3,5-Trimethylbenzene	9.7	10.0	97.0%	9.7	10.0	97.0% 0.0%
1,2,4-Trimethylbenzene	9.7	10.0	97.0%	9.9	10.0	99.0% 2.0%
Hexachlorobutadiene	9.8	10.0	98.0%	10.3	10.0	103% 5.0%
Ethylene Dibromide	9.6	10.0	96.0%	9.9	10.0	99.0% 3.1%
Bromochloromethane	9.7	10.0	97.0%	9.8	10.0	98.0% 1.0%
2,2-Dichloropropane	9.0	10.0	90.0%	9.0	10.0	90.0% 0.0%
1,3-Dichloropropane	9.5	10.0	95.0%	9.5	10.0	95.0% 0.0%
Isopropylbenzene	9.7	10.0	97.0%	9.7	10.0	97.0% 0.0%
n-Propylbenzene	9.5	10.0	95.0%	9.7	10.0	97.0% 2.1%
Bromobenzene	9.7	10.0	97.0%	9.7	10.0	97.0% 0.0%
2-Chlorotoluene	9.4	10.0	94.0%	9.5	10.0	95.0% 1.1%
4-Chlorotoluene	9.5	10.0	95.0%	9.6	10.0	96.0% 1.0%
tert-Butylbenzene	9.6	10.0	96.0%	9.7	10.0	97.0% 1.0%
sec-Butylbenzene	9.5	10.0	95.0%	9.8	10.0	98.0% 3.1%
4-Isopropyltoluene	9.6	10.0	96.0%	10.0	10.0	100% 4.1%
n-Butylbenzene	9.0	10.0	90.0%	9.7	10.0	97.0% 7.5%
1,2,4-Trichlorobenzene	9.8	10.0	98.0%	9.9	10.0	99.0% 1.0%
Naphthalene	10.2	10.0	102%	10.3	10.0	103% 1.0%
1,2,3-Trichlorobenzene	10.0	10.0	100%	10.2	10.0	102% 2.0%

Reported in µg/L (ppb)

RPD calculated using sample concentrations per SW846.

Volatile Surrogate Recovery

	LCS	LCSD
d4-1,2-Dichloroethane	93.9%	93.3%
d8-Toluene	98.3%	99.9%
Bromofluorobenzene	98.1%	98.3%
d4-1,2-Dichlorobenzene	101%	99.9%

4A
VOLATILE METHOD BLANK SUMMARY

Method Blank ID.

	MB0926
--	--------

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONME

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVE

Lab File ID: MB0926

Lab Sample ID: MB0926

Date Analyzed: 09/26/11

Time Analyzed: 1100

Instrument ID: NT2

Heated Purge: (Y/N) N

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01 LCS0926	LCS0926	LCS0926	1006
02 LCS0926	LCS0926	LCS0926A	1033
03 TRIP BLANKS	TN42G	TN42G2	1228
04 FRP-091911-0	TN00B	TN00B2	1255
05 FRP-091911-0	TN00C	TN00C2	1322
06 FRP-091911-0	TN00D	TN00D2	1349
07 FRP-091911-0	TN00E	TN00E2	1415
08 FRP-092011-0	TN19B	TN19B2	1442
09 FRP-092011-0	TN19C	TN19C2	1509
10 FRP-092011-0	TN19D	TN19D2	1536
11 FRP-092111-0	TN42A	TN42A2	1603
12 FRP-092111-0	TN42B	TN42B2	1630
13 FRP-092111-0	TN42C	TN42C2	1657
14 FRP-092111-0	TN42D	TN42D2	1723
15 FRP-092111-0	TN42E	TN42E2	1750
16 FRP-092111-0	TN42F	TN42F2	1817
17 FRP-091911-0	TN00CMS	TN00CMS	1843
18 FRP-091911-0	TN00CMSP	TN00CMSP	1910
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COMMENTS:

ORGANICS ANALYSIS DATA SHEET

Volatile s by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2

**ANALYTICAL
RESOURCES
INCORPORATED**

Sample ID: MB-092611

METHOD BLANK

Lab Sample ID: MB-092611

LIMS ID: 11-20663

Matrix: Water

Data Release Authorized: MW

Reported: 10/03/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: NA

Date Received: NA

Instrument/Analyst: NT2/PAB

Date Analyzed: 09/26/11 11:00

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

Sample ID: MB-092611
METHOD BLANK

Lab Sample ID: MB-092611
LIMS ID: 11-20663
Matrix: Water
Date Analyzed: 09/26/11 11:00

QC Report No: TN42-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromochloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	94.0%
d8-Toluene	99.0%
Bromofluorobenzene	96.0%
d4-1,2-Dichlorobenzene	100%

5A
VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: ANALYTICAL RESOURCES INC Contract: AMEC EARTH AND ENVIRONMENTAL
 Lab Code: ARI Case No.: 2011 FRP SHORELINE INVESTIGATION SDG No.: TN42
 Lab File ID: BFB0727 BFB Injection Date: 07/27/11
 Instrument ID: NT5 BFB Injection Time: 1130
 GC Column: RTXVMS ID: 0.18 (mm) Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	8.0 - 40.0% of mass 95	16.1
75	30.0 - 66.0% of mass 95	55.5
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.0
173	Less than 2.0% of mass 174	0.7 (1.0)1
174	50.0 - 101.0% of mass 95	72.0
175	4.0 - 9.0% of mass 174	6.0 (8.4)1
176	93.0 - 101.0% of mass 174	72.1 (100.1)1
177	5.0 - 9.0% of mass 176	4.7 (6.5)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01 VSTD60	IC0727	6000727	07/27/11	1220
02 VSTD40	IC0727	4000727	07/27/11	1248
03 VSTD20	IC0727	2000727	07/27/11	1316
04 VSTD10	IC0727	1000727	07/27/11	1345
05 VSTD4	IC0727	0400727	07/27/11	1413
06 VSTD1	IC0727	0100727	07/27/11	1441
07 VSTD0.5	IC0727	0050727	07/27/11	1510
08 VSTD0.2	IC0727	0020727	07/27/11	1538
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5A
 VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
 BROMOFLUOROBENZENE (BFB)

Lab Name: ANALYTICAL RESOURCES INC Contract: AMEC EARTH AND ENVIRONMENTAL
 Lab Code: ARI Case No.: 2011 FRP SHORELINE INVESTIGATION SDG No.: TN42
 Lab File ID: BFB0921 BFB Injection Date: 09/21/11
 Instrument ID: NT5 BFB Injection Time: 0812
 GC Column: RTXVMS ID: 0.18 (mm) Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	8.0 - 40.0% of mass 95	18.7
75	30.0 - 66.0% of mass 95	55.0
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	5.9
173	Less than 2.0% of mass 174	0.2 (0.2)1
174	50.0 - 101.0% of mass 95	73.7
175	4.0 - 9.0% of mass 174	5.4 (7.3)1
176	93.0 - 101.0% of mass 174	73.3 (99.4)1
177	5.0 - 9.0% of mass 176	4.4 (5.9)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01 VSTD10	CC0921	1000921	09/21/11	0855
02 LCS0921	LCS0921	LCS0921	09/21/11	1001
03 LCS0921	LCS0921	LCS0921A	09/21/11	1029
04 MB0921	MB0921	MB0921	09/21/11	1057
05 FRP-091911-002	TN00B	TN00B	09/21/11	1433
06 FRP-091911-003	TN00C	TN00C	09/21/11	1501
07 FRP-091911-004	TN00D	TN00D	09/21/11	1529
08 FRP-091911-005	TN00E	TN00E	09/21/11	1557
09 FRP-091911-006	TN00F	TN00F	09/21/11	1626
10 TRIP BLANKS	TN00G	TN00G	09/21/11	1654
11 FRP-092011-001	TN19A	TN19A	09/21/11	1722
12 FRP-092011-005	TN19E	TN19E	09/21/11	1915
13 TRIP BLANKS	TN19F	TN19F	09/21/11	1944
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5A
VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: ANALYTICAL RESOURCES INC Contract: AMEC EARTH AND ENVIRONMENTAL
 Lab Code: ARI Case No.: 2011 FRP SHORELINE INVESTIGATION SDG No.: TN42
 Lab File ID: BFB0921 BFB Injection Date: 09/21/11
 Instrument ID: NT2 BFB Injection Time: 1241
 GC Column: RTXVMS ID: 0.18 (mm) Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	8.0 - 40.0% of mass 95	18.7
75	30.0 - 66.0% of mass 95	53.0
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.1
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	50.0 - 101.0% of mass 95	79.2
175	4.0 - 9.0% of mass 174	5.7 (7.2)1
176	93.0 - 101.0% of mass 174	79.4 (100.3)1
177	5.0 - 9.0% of mass 176	5.5 (7.0)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01 VSTD0.2	VSTD0.2	00_20921	09/21/11	1338
02 VSTD0.5	VSTD0.5	00_50921	09/21/11	1405
03 VSTD01	VSTD01	01_00921	09/21/11	1432
04 VSTD02	VSTD02	02_00921	09/21/11	1459
05 VSTD10	VSTD10	10_00921	09/21/11	1526
06 VSTD20	VSTD20	20_00921	09/21/11	1552
07 VSTD40	VSTD40	40_00921	09/21/11	1619
08 VSTD60	VSTD60	60_00921	09/21/11	1646
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5A
VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: ANALYTICAL RESOURCES INC Contract: AMEC EARTH AND ENVIRONMENTAL
 Lab Code: ARI Case No.: 2011 FRP SHORELINE INVESTIGATION SDG No.: TN42
 Lab File ID: BFB0926 BFB Injection Date: 09/26/11
 Instrument ID: NT2 BFB Injection Time: 0901
 GC Column: RTXVMS ID: 0.18 (mm) Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	8.0 - 40.0% of mass 95	17.4
75	30.0 - 66.0% of mass 95	51.2
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.2
173	Less than 2.0% of mass 174	0.1 (0.1)1
174	50.0 - 101.0% of mass 95	82.2
175	4.0 - 9.0% of mass 174	5.7 (6.9)1
176	93.0 - 101.0% of mass 174	81.9 (99.6)1
177	5.0 - 9.0% of mass 176	5.5 (6.7)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	CC0926	CC0926	CC0926	09/26/11	0939
02	LCS0926	LCS0926	LCS0926	09/26/11	1006
03	LCS0926	LCS0926	LCS0926A	09/26/11	1033
04	MB0926	MB0926	MB0926	09/26/11	1100
05	TRIP BLANKS	TN42G	TN42G2	09/26/11	1228
06	FRP-091911-002	TN00B	TN00B2	09/26/11	1255
07	FRP-091911-003	TN00C	TN00C2	09/26/11	1322
08	FRP-091911-004	TN00D	TN00D2	09/26/11	1349
09	FRP-091911-005	TN00E	TN00E2	09/26/11	1415
10	FRP-092011-002	TN19B	TN19B2	09/26/11	1442
11	FRP-092011-003	TN19C	TN19C2	09/26/11	1509
12	FRP-092011-004	TN19D	TN19D2	09/26/11	1536
13	FRP-092111-001	TN42A	TN42A2	09/26/11	1603
14	FRP-092111-002	TN42B	TN42B2	09/26/11	1630
15	FRP-092111-003	TN42C	TN42C2	09/26/11	1657
16	FRP-092111-004	TN42D	TN42D2	09/26/11	1723
17	FRP-092111-005	TN42E	TN42E2	09/26/11	1750
18	FRP-092111-006	TN42F	TN42F2	09/26/11	1817
19	FRP-091911-003MS	TN00CMS	TN00CMS	09/26/11	1843
20	FRP-091911-003MS	TN00CMSP	TN00CMSP	09/26/11	1910
21					
22					

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF0.2: 0020727 RF0.5: 0050727 RF1: 0100727
 RF4: 0400727 RF10: 1000727

COMPOUND	RF0.2	RF0.5	RF1	RF4	RF10
Chloromethane		0.445	0.457	0.360	0.340
Vinyl Chloride	0.387	0.360	0.422	0.419	0.431
Bromomethane	0.197	0.224	0.234	0.250	0.262
Chloroethane	0.276	0.227	0.280	0.231	0.255
Trichlorofluoromethane	0.957	0.998	1.258	1.074	1.068
Acrolein		0.042	0.034	0.033	0.034
112Trichloro122Trifluoroetha	0.463	0.450	0.550	0.466	0.445
Acetone			0.061	0.057	0.063
1,1-Dichloroethene	0.352	0.371	0.490	0.397	0.396
Bromoethane	0.262	0.252	0.278	0.248	0.257
Iodomethane		0.564	0.749	0.624	0.606
Methylene Chloride	0.348	0.486	0.452	0.387	0.396
Acrylonitrile			0.089	0.069	0.060
Carbon Disulfide	1.156	1.041	1.184	1.061	1.118
Trans-1,2-Dichloroethene	0.430	0.490	0.510	0.467	0.460
Vinyl Acetate		0.413	0.474	0.465	0.461
1,1-Dichloroethane	0.679	0.751	0.805	0.733	0.736
2-Butanone		0.078	0.097	0.087	0.083
2,2-Dichloropropane	0.862	0.880	1.027	0.962	0.990
Cis-1,2-Dichloroethene	0.446	0.434	0.510	0.463	0.474
Chloroform	1.037	0.831	1.051	0.976	0.911
Bromochloromethane	0.185	0.222	0.241	0.232	0.214
1,1,1-Trichloroethane	0.935	1.018	1.168	1.118	1.080
1,1-Dichloropropene	0.495	0.433	0.524	0.489	0.504
Carbon Tetrachloride	0.648	0.697	0.815	0.779	0.777
1,2-Dichloroethane	0.571	0.510	0.641	0.592	0.574
Benzene	1.219	1.141	1.347	1.264	1.263
Trichloroethene	0.359	0.361	0.405	0.402	0.397
1,2-Dichloropropane	0.197	0.242	0.301	0.233	0.240
Bromodichloromethane	0.587	0.499	0.577	0.564	0.540
Dibromomethane	0.172	0.180	0.213	0.200	0.194
2-Chloroethyl Vinyl Ether		0.094	0.116	0.115	0.124
4-Methyl-2-Pentanone		0.058	0.064	0.063	0.062
Cis 1,3-dichloropropene	0.466	0.524	0.587	0.555	0.548
Toluene	0.906	0.806	0.927	0.902	0.904
Trans 1,3-Dichloropropene	0.529	0.478	0.594	0.558	0.556
2-Hexanone		0.099	0.124	0.114	0.115

FORM VI VOA

TN42 : 022945

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF0.2: 0020727 RF0.5: 0050727 RF1: 0100727
 RF4: 0400727 RF10: 1000727

COMPOUND	RF0.2	RF0.5	RF1	RF4	RF10
1,1,2-Trichloroethane	0.227	0.237	0.281	0.249	0.239
1,3-Dichloropropane	0.471	0.429	0.523	0.470	0.439
Tetrachloroethene	0.510	0.433	0.513	0.467	0.454
Chlorodibromomethane	0.381	0.358	0.464	0.425	0.420
1,2-Dibromoethane	0.280	0.234	0.279	0.264	0.262
Chlorobenzene	1.158	1.115	1.204	1.137	1.105
Ethyl Benzene	0.584	0.562	0.651	0.606	0.613
1,1,1,2-Tetrachloroethane	0.450	0.454	0.514	0.471	0.461
m,p-xylene	0.719	0.700	0.823	0.767	0.747
o-Xylene	0.592	0.609	0.769	0.695	0.714
Styrene	0.995	1.042	1.262	1.183	1.210
Bromoform	0.486	0.394	0.515	0.437	0.411
1,1,2,2-Tetrachloroethane	0.675	0.420	0.492	0.469	0.413
1,2,3-Trichloropropane	0.264	0.209	0.283	0.207	0.191
Trans-1,4-Dichloro 2-Butene			0.229	0.236	0.207
N-Propyl Benzene	3.544	3.127	3.884	3.502	3.307
Bromobenzene	0.950	0.906	1.048	0.900	0.835
Isopropyl Benzene	2.811	2.776	3.485	3.289	3.092
2-Chloro Toluene	2.351	2.305	2.711	2.443	2.307
4-Chloro Toluene	2.961	2.400	2.912	2.656	2.526
T-Butyl Benzene	2.160	2.101	2.624	2.386	2.279
1,3,5-Trimethyl Benzene	2.657	2.500	3.198	2.913	2.747
1,2,4-Trimethylbenzene	2.740	2.526	3.143	2.980	2.852
S-Butyl Benzene	3.035	2.770	3.343	2.996	2.938
4-Isopropyl Toluene	2.488	2.252	2.859	2.751	2.701
1,3-Dichlorobenzene	1.678	1.489	1.999	1.701	1.598
1,4-Dichlorobenzene	2.105	1.802	1.971	1.734	1.649
N-Butyl Benzene	1.928	1.951	2.329	2.148	2.111
1,2-Dichlorobenzene	1.733	1.498	1.780	1.579	1.456
1,2-Dibromo 3-Chloropropane	0.104	0.177	0.173	0.133	0.120
1,2,4-Trichlorobenzene	0.910	0.873	1.027	0.890	0.833
Hexachloro 1,3-Butadiene		0.550	0.577	0.472	0.444
Naphthalene	1.426	1.383	1.643	1.510	1.454
1,2,3-Trichlorobenzene	0.750	0.612	0.758	0.638	0.546
Dichlorodifluoromethane	0.511	0.578	0.594	0.616	0.646
Methyl tert butyl ether	1.258	1.246	1.428	1.319	1.291

FORM VI VOA

TN42 : 000060

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF0.2: 0020727 RF0.5: 0050727 RF1: 0100727
RF4: 0400727 RF10: 1000727

COMPOUND	RF0.2	RF0.5	RF1	RF4	RF10
d4-1,2-Dichloroethane	0.541	0.588	0.595	0.598	0.570
d8-Toluene	1.120	1.100	1.111	1.136	1.126
4-Bromofluorobenzene	0.554	0.579	0.575	0.589	0.596
d4-1,2-Dichlorobenzene	0.861	0.875	0.902	0.882	0.882
Dibromofluoromethane	0.399	0.403	0.423	0.423	0.426

FORM VI VOA

TN42 : 000851

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF20: 2000727

RF40: 4000727

RF60: 6000727

COMPOUND	RF20	RF40	RF60
Chloromethane	0.307	0.274	0.259
Vinyl Chloride	0.382	0.362	0.311
Bromomethane	0.267	0.276	0.256
Chloroethane	0.215	0.210	0.203
Trichlorofluoromethane	0.942	0.762	0.719
Acrolein	0.031	0.029	0.028
112Trichloro122Trifluoroetha	0.409	0.391	0.374
Acetone	0.049	0.054	0.053
1,1-Dichloroethene	0.359	0.339	0.314
Bromoethane	0.233	0.207	0.180
Iodomethane	0.548	0.505	0.418
Methylene Chloride	0.356	0.322	0.283
Acrylonitrile	0.060	0.059	0.058
Carbon Disulfide	0.995	0.897	0.942
Trans-1,2-Dichloroethene	0.437	0.405	0.414
Vinyl Acetate	0.456	0.463	0.423
1,1-Dichloroethane	0.690	0.698	0.621
2-Butanone	0.074	0.077	0.070
2,2-Dichloropropane	0.931	0.929	0.854
Cis-1,2-Dichloroethene	0.436	0.435	0.401
Chloroform	0.855	0.847	0.762
Bromochloromethane	0.200	0.194	0.182
1,1,1-Trichloroethane	0.993	0.975	0.872
1,1-Dichloropropene	0.468	0.452	0.424
Carbon Tetrachloride	0.718	0.701	0.637
1,2-Dichloroethane	0.540	0.532	0.492
Benzene	1.178	1.121	1.043
Trichloroethene	0.371	0.360	0.331
1,2-Dichloropropane	0.229	0.218	0.202
Bromodichloromethane	0.497	0.481	0.447
Dibromomethane	0.176	0.169	0.158
2-Chloroethyl Vinyl Ether	0.115	0.113	0.109
4-Methyl-2-Pentanone	0.057	0.054	0.050
Cis 1,3-dichloropropene	0.503	0.490	0.458
Toluene	0.830	0.800	0.734
Trans 1,3-Dichloropropene	0.525	0.499	0.455
2-Hexanone	0.102	0.080	

FORM VI VOA

TN42 : 00052

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF20: 2000727

RF40: 4000727

RF60: 6000727

COMPOUND	RF20	RF40	RF60
1,1,2-Trichloroethane	0.221	0.218	0.204
1,3-Dichloropropane	0.402	0.396	0.376
Tetrachloroethene	0.430	0.409	0.384
Chlorodibromomethane	0.378	0.371	0.347
1,2-Dibromoethane	0.244	0.230	0.219
Chlorobenzene	1.021	0.986	0.894
Ethyl Benzene	0.575	0.540	0.497
1,1,1,2-Tetrachloroethane	0.426	0.411	0.376
m,p-xylene	0.690	0.647	0.578
o-Xylene	0.676	0.648	0.591
Styrene	1.125	1.045	0.944
Bromoform	0.358	0.339	
1,1,2,2-Tetrachloroethane	0.405		
1,2,3-Trichloropropane	0.174		
Trans-1,4-Dichloro 2-Butene	0.184	0.132	
N-Propyl Benzene	3.004	2.929	2.640
Bromobenzene	0.746	0.727	0.648
Isopropyl Benzene	2.818	2.721	2.425
2-Chloro Toluene	2.102	2.003	1.874
4-Chloro Toluene	2.340	2.147	2.131
T-Butyl Benzene	2.101	2.066	1.893
1,3,5-Trimethyl Benzene	2.529	2.467	2.227
1,2,4-Trimethylbenzene	2.608	2.552	2.326
S-Butyl Benzene	2.733	2.654	2.408
4-Isopropyl Toluene	2.494	2.465	2.192
1,3-Dichlorobenzene	1.470		
1,4-Dichlorobenzene	1.506	1.492	1.294
N-Butyl Benzene	1.986	1.974	
1,2-Dichlorobenzene	1.375	1.334	1.156
1,2-Dibromo 3-Chloropropane	0.113	0.105	0.090
1,2,4-Trichlorobenzene	0.766		
Hexachloro 1,3-Butadiene	0.385		
Naphthalene	1.353		
1,2,3-Trichlorobenzene	0.398		
Dichlorodifluoromethane	0.604	0.542	0.459
Methyl tert butyl ether	1.204	1.215	1.114

FORM VI VOA

TN42:000053

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF20: 2000727 RF40: 4000727 RF60: 6000727

COMPOUND	RF20	RF40	RF60
d4-1,2-Dichloroethane	0.570	0.588	0.563
d8-Toluene	1.125	1.115	1.108
4-Bromofluorobenzene	0.620	0.597	0.505
d4-1,2-Dichlorobenzene	0.894	0.880	0.839
Dibromofluoromethane	0.415	0.422	0.409

FORM VI VOA

TN42 : 000054

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R ²
Chloromethane	LINR		0.9926
Vinyl Chloride	AVRG	0.384	10.4
Bromomethane	AVRG	0.246	10.5
Chloroethane	AVRG	0.237	12.5
Trichlorofluoromethane	AVRG	0.972	17.8
Acrolein	AVRG	0.033	13.6
112Trichloro122Trifluoroetha	AVRG	0.444	12.3
Acetone	AVRG	0.056	9.7
1,1-Dichloroethene	AVRG	0.377	14.1
Bromoethane	AVRG	0.240	13.4
Iodomethane	AVRG	0.573	18.0
Methylene Chloride	AVRG	0.379	17.6
Acrylonitrile	AVRG	0.066	18.2
Carbon Disulfide	AVRG	1.049	9.7
Trans-1,2-Dichloroethene	AVRG	0.452	8.1
Vinyl Acetate	AVRG	0.451	5.1
1,1-Dichloroethane	AVRG	0.714	7.7
2-Butanone	AVRG	0.081	11.3
2,2-Dichloropropane	AVRG	0.929	6.7
Cis-1,2-Dichloroethene	AVRG	0.450	7.3
Chloroform	AVRG	0.909	11.4
Bromochloromethane	AVRG	0.209	10.5
1,1,1-Trichloroethane	AVRG	1.020	9.6
1,1-Dichloropropene	AVRG	0.474	7.4
Carbon Tetrachloride	AVRG	0.722	8.9
1,2-Dichloroethane	AVRG	0.556	8.6
Benzene	AVRG	1.197	8.0
Trichloroethene	AVRG	0.373	7.0
1,2-Dichloropropane	AVRG	0.233	13.8
Bromodichloromethane	AVRG	0.524	9.6
Dibromomethane	AVRG	0.183	9.9
2-Chloroethyl Vinyl Ether	AVRG	0.112	8.4
4-Methyl-2-Pentanone	AVRG	0.058	8.9
Cis 1,3-dichloropropene	AVRG	0.516	8.8
Toluene	AVRG	0.851	8.1
Trans 1,3-Dichloropropene	AVRG	0.524	8.7
2-Hexanone	AVRG	0.105	14.6

<- Indicates value outside QC limits:
(%RSD < 20% or R² > 0.990)

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R ²
1,1,2-Trichloroethane	AVRG	0.234	10.0
1,3-Dichloropropane	AVRG	0.438	11.0
Tetrachloroethene	AVRG	0.450	10.1
Chlorodibromomethane	AVRG	0.393	10.1
1,2-Dibromoethane	AVRG	0.252	9.2
Chlorobenzene	AVRG	1.078	9.5
Ethyl Benzene	AVRG	0.578	8.2
1,1,1,2-Tetrachloroethane	AVRG	0.446	9.3
m,p-xylene	AVRG	0.709	10.6
o-Xylene	AVRG	0.662	9.6
Styrene	AVRG	1.101	10.2
Bromoform	AVRG	0.420	15.3
1,1,2,2-Tetrachloroethane	LINR		0.9985
1,2,3-Trichloropropane	AVRG	0.221	19.2
Trans-1,4-Dichloro 2-Butene	2ORDR		0.9956
N-Propyl Benzene	AVRG	3.242	12.2
Bromobenzene	AVRG	0.845	15.6
Isopropyl Benzene	AVRG	2.927	11.6
2-Chloro Toluene	AVRG	2.262	11.7
4-Chloro Toluene	AVRG	2.509	12.6
T-Butyl Benzene	AVRG	2.201	10.2
1,3,5-Trimethyl Benzene	AVRG	2.655	11.3
1,2,4-Trimethylbenzene	AVRG	2.716	9.8
S-Butyl Benzene	AVRG	2.860	9.9
4-Isopropyl Toluene	AVRG	2.525	9.3
1,3-Dichlorobenzene	AVRG	1.656	11.6
1,4-Dichlorobenzene	AVRG	1.694	15.7
N-Butyl Benzene	AVRG	2.061	7.0
1,2-Dichlorobenzene	AVRG	1.489	13.9
1,2-Dibromo 3-Chloropropane	2ORDR		0.9984
1,2,4-Trichlorobenzene	AVRG	0.883	9.8
Hexachloro 1,3-Butadiene	AVRG	0.485	16.1
Naphthalene	AVRG	1.461	7.1
1,2,3-Trichlorobenzene	2ORDR		0.9972
Dichlorodifluoromethane	AVRG	0.569	10.7
Methyl tert butyl ether	AVRG	1.260	7.3

<- Indicates value outside QC limits:
(%RSD < 20% or R² > 0.990)

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R ²
d4-1,2-Dichloroethane	AVRG	0.577	3.4
d8-Toluene	AVRG	1.118	1.0
4-Bromofluorobenzene	AVRG	0.577	6.0
d4-1,2-Dichlorobenzene	AVRG	0.877	2.2
Dibromofluoromethane	AVRG	0.415	2.5

<- Indicates value outside QC limits:
(%RSD < 20% or R² > 0.990)

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF0.2: 00_20921 RF0.5: 00_50921 RF1: 01_00921
 RF2: 02_00921 RF10: 10_00921

COMPOUND	RF0.2	RF0.5	RF1	RF2	RF10
Chloromethane		0.942	0.850	0.880	0.733
Vinyl Chloride	0.625	0.610	0.625	0.659	0.642
Bromomethane			0.323	0.325	0.307
Chloroethane	0.427	0.407	0.388	0.412	0.385
Trichlorofluoromethane	0.798	0.711	0.725	0.783	0.724
Acrolein	0.058	0.054	0.057	0.057	0.061
112Trichloro122Trifluoroetha	0.529	0.449	0.475	0.479	0.440
Acetone		0.231	0.193	0.173	0.148
1,1-Dichloroethene	0.625	0.508	0.473	0.498	0.444
Bromoethane	0.340	0.311	0.329	0.338	0.314
Iodomethane			0.526	0.521	0.551
Methylene Chloride		0.598	0.546	0.550	0.507
Acrylonitrile		0.155	0.158	0.169	0.160
Carbon Disulfide	1.638	1.529	1.519	1.651	1.514
Trans-1,2-Dichloroethene	0.564	0.536	0.523	0.543	0.502
Vinyl Acetate	0.978	0.950	0.997	1.058	1.040
1,1-Dichloroethane	1.007	0.928	0.954	0.988	0.910
2-Butanone	0.252	0.236	0.247	0.257	0.244
2,2-Dichloropropane	0.928	0.805	0.770	0.815	0.778
Cis-1,2-Dichloroethene	0.584	0.561	0.559	0.573	0.534
Chloroform	0.956	0.887	0.938	0.963	0.909
Bromochloromethane	0.265	0.250	0.254	0.272	0.250
1,1,1-Trichloroethane	0.897	0.862	0.869	0.900	0.842
1,1-Dichloropropene	0.509	0.504	0.492	0.522	0.490
Carbon Tetrachloride	0.456	0.431	0.439	0.462	0.459
1,2-Dichloroethane	0.498	0.505	0.489	0.510	0.455
Benzene	1.487	1.436	1.433	1.489	1.356
Trichloroethene	0.407	0.350	0.365	0.370	0.346
1,2-Dichloropropane	0.383	0.348	0.350	0.373	0.339
Bromodichloromethane	0.450	0.418	0.421	0.454	0.447
Dibromomethane	0.205	0.183	0.200	0.207	0.187
2-Chloroethyl Vinyl Ether		0.117	0.114	0.107	0.146
4-Methyl-2-Pentanone	0.118	0.126	0.131	0.140	0.135
Cis 1,3-dichloropropene	0.547	0.501	0.535	0.563	0.552
Toluene	0.946	0.932	0.930	0.961	0.874
Trans 1,3-Dichloropropene	0.468	0.467	0.480	0.522	0.521
2-Hexanone	0.251	0.252	0.262	0.286	0.280

FORM VI VOA

TN42:000050

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF0.2: 00_20921 RF0.5: 00_50921 RF1: 01_00921
 RF2: 02_00921 RF10: 10_00921

COMPOUND	RF0.2	RF0.5	RF1	RF2	RF10
1,1,2-Trichloroethane	0.328	0.290	0.297	0.308	0.286
1,3-Dichloropropane	0.575	0.562	0.574	0.590	0.558
Tetrachloroethene	0.412	0.403	0.418	0.415	0.394
Chlorodibromomethane	0.263	0.289	0.292	0.317	0.329
1,2-Dibromoethane	0.286	0.288	0.286	0.320	0.301
Chlorobenzene	1.135	1.029	1.070	1.096	1.005
Ethyl Benzene	0.608	0.561	0.552	0.588	0.548
1,1,1,2-Tetrachloroethane	0.347	0.304	0.329	0.358	0.353
m,p-xylene	0.689	0.688	0.693	0.737	0.677
o-Xylene	0.662	0.663	0.679	0.716	0.659
Styrene	0.991	1.031	1.056	1.187	1.121
Bromoform	0.280	0.276	0.289	0.304	0.343
1,1,2,2-Tetrachloroethane	0.602	0.660	0.667	0.683	0.684
1,2,3-Trichloropropane	0.250	0.226	0.232	0.229	0.226
Trans-1,4-Dichloro 2-Butene			0.119	0.136	0.125
N-Propyl Benzene	3.640	3.514	3.562	3.710	3.528
Bromobenzene	0.829	0.749	0.778	0.784	0.746
Isopropyl Benzene	2.501	2.530	2.531	2.670	2.579
2-Chloro Toluene	2.555	2.513	2.509	2.536	2.393
4-Chloro Toluene	2.419	2.278	2.373	2.345	2.236
T-Butyl Benzene	2.417	2.330	2.302	2.437	2.316
1,3,5-Trimethyl Benzene	2.501	2.530	2.531	2.670	2.579
1,2,4-Trimethylbenzene	2.527	2.071	2.293	2.468	2.448
S-Butyl Benzene	3.444	3.038	3.114	3.397	3.276
4-Isopropyl Toluene	2.861	2.170	2.452	2.708	2.665
1,3-Dichlorobenzene	1.651	1.547	1.536	1.610	1.506
1,4-Dichlorobenzene	1.674	1.654	1.641	1.681	1.556
N-Butyl Benzene	2.395	1.518	1.725	2.134	2.087
1,2-Dichlorobenzene	1.615	1.556	1.570	1.588	1.471
1,2-Dibromo 3-Chloropropane		0.112	0.127	0.133	0.135
1,2,4-Trichlorobenzene	1.082	0.757	0.815	0.997	0.996
Hexachloro 1,3-Butadiene		0.559	0.560	0.600	0.576
Naphthalene		1.645	1.739	2.138	2.152
1,2,3-Trichlorobenzene		0.704	0.719	0.921	0.886
Dichlorodifluoromethane	0.428	0.483	0.425	0.502	0.483
Methyl tert butyl ether	1.458	1.374	1.459	1.378	1.400

FORM VI VOA

TN42:00059

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF0.2: 00_20921 RF0.5: 00_50921 RF1: 01_00921
RF2: 02_00921 RF10: 10_00921

COMPOUND	RF0.2	RF0.5	RF1	RF2	RF10
d4-1,2-Dichloroethane	0.582	0.575	0.582	0.580	0.575
d8-Toluene	1.240	1.232	1.232	1.241	1.212
4-Bromofluorobenzene	0.558	0.551	0.550	0.557	0.548
d4-1,2-Dichlorobenzene	0.906	0.898	0.914	0.904	0.910
Dibromofluoromethane	0.449	0.444	0.447	0.453	0.459

FORM VI VOA

TN42:00066

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF20: 20_00921 RF40: 40_00921 RF60: 60_00921

COMPOUND	RF20	RF40	RF60
Chloromethane	0.729	0.758	0.786
Vinyl Chloride	0.640	0.684	0.645
Bromomethane	0.311	0.341	0.355
Chloroethane	0.329		
Trichlorofluoromethane	0.725	0.682	0.628
Acrolein	0.070	0.071	0.073
112Trichloro122Trifluoroetha	0.460	0.467	0.457
Acetone	0.156	0.155	0.157
1,1-Dichloroethene	0.457	0.462	0.460
Bromoethane	0.316	0.313	0.311
Iodomethane	0.545	0.517	0.534
Methylene Chloride	0.524	0.536	0.534
Acrylonitrile	0.172	0.176	0.178
Carbon Disulfide	1.611	1.623	1.614
Trans-1,2-Dichloroethene	0.526	0.528	0.532
Vinyl Acetate	1.115	1.160	1.165
1,1-Dichloroethane	0.954	0.972	0.980
2-Butanone	0.265	0.270	
2,2-Dichloropropane	0.814	0.826	0.822
Cis-1,2-Dichloroethene	0.558	0.567	0.565
Chloroform	0.952	0.968	0.972
Bromochloromethane	0.258	0.259	0.261
1,1,1-Trichloroethane	0.890	0.900	0.893
1,1-Dichloropropene	0.517	0.519	0.517
Carbon Tetrachloride	0.492	0.499	0.506
1,2-Dichloroethane	0.486	0.485	0.492
Benzene	1.419	1.402	1.387
Trichloroethene	0.362	0.363	0.366
1,2-Dichloropropane	0.358	0.361	0.364
Bromodichloromethane	0.488	0.492	0.500
Dibromomethane	0.199	0.198	0.199
2-Chloroethyl Vinyl Ether	0.158	0.162	0.173
4-Methyl-2-Pentanone	0.147	0.144	0.147
Cis 1,3-dichloropropene	0.604	0.610	0.616
Toluene	0.914	0.897	0.892
Trans 1,3-Dichloropropene	0.566	0.576	0.582
2-Hexanone	0.296	0.290	0.279

FORM VI VOA

TN42 : 000961

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF20: 20_00921 RF40: 40_00921 RF60: 60_00921

COMPOUND	RF20	RF40	RF60
1,1,2-Trichloroethane	0.304	0.301	0.304
1,3-Dichloropropane	0.575	0.587	0.583
Tetrachloroethene	0.409	0.410	0.404
Chlorodibromomethane	0.360	0.377	0.379
1,2-Dibromoethane	0.318	0.320	0.326
Chlorobenzene	1.037	1.035	1.023
Ethyl Benzene	0.569	0.573	0.571
1,1,1,2-Tetrachloroethane	0.370	0.380	0.382
m,p-xylene	0.700	0.691	0.678
c-Xylene	0.682	0.690	0.695
Styrene	1.186	1.225	1.221
Bromoform	0.409	0.422	0.429
1,1,2,2-Tetrachloroethane	0.728	0.727	0.728
1,2,3-Trichloropropane	0.239	0.234	0.231
Trans-1,4-Dichloro 2-Butene	0.178	0.194	0.200
N-Propyl Benzene	3.624	3.444	3.194
Bromobenzene	0.780	0.778	0.774
Isopropyl Benzene	2.668	2.616	2.475
2-Chloro Toluene	2.494	2.461	2.388
4-Chloro Toluene	2.320	2.283	2.204
T-Butyl Benzene	2.394	2.368	2.271
1,3,5-Trimethyl Benzene	2.668	2.616	2.475
1,2,4-Trimethylbenzene	2.601	2.622	2.505
S-Butyl Benzene	3.306	3.277	3.045
4-Isopropyl Toluene	2.793	2.859	2.682
1,3-Dichlorobenzene	1.565	1.555	1.525
1,4-Dichlorobenzene	1.619	1.606	1.569
N-Butyl Benzene	2.079	2.322	2.111
1,2-Dichlorobenzene	1.524	1.514	1.485
1,2-Dibromo 3-Chloropropane	0.153	0.155	0.158
1,2,4-Trichlorobenzene	0.955	1.014	0.957
Hexachloro 1,3-Butadiene	0.572	0.560	0.548
Naphthalene	1.962	1.832	1.857
1,2,3-Trichlorobenzene	0.865	0.888	0.848
Dichlorodifluoromethane	0.488	0.513	0.472
Methyl tert butyl ether	1.496	1.501	1.544

FORM VI VOA

TN42 : 000062

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF20: 20_00921 RF40: 40_00921 RF60: 60_00921

COMPOUND	RF20	RF40	RF60
d4-1,2-Dichloroethane	0.581	0.588	0.588
d8-Toluene	1.242	1.215	1.221
4-Bromofluorobenzene	0.550	0.558	0.573
d4-1,2-Dichlorobenzene	0.918	0.909	0.903
Dibromofluoromethane	0.461	0.470	0.478

FORM VI VOA

TN42 : 000063

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R ²
Chloromethane	AVRG	0.811	10.0
Vinyl Chloride	AVRG	0.641	3.5
Bromomethane	AVRG	0.327	5.6
Chloroethane	AVRG	0.391	8.7
Trichlorofluoromethane	AVRG	0.722	7.4
Acrolein	AVRG	0.063	11.9
112Trichloro122Trifluoroetha	AVRG	0.470	5.8
Acetone	AVRG	0.173	17.2
1,1-Dichloroethene	AVRG	0.491	11.9
Bromoethane	AVRG	0.321	3.8
Iodomethane	AVRG	0.532	2.6
Methylene Chloride	AVRG	0.542	5.3
Acrylonitrile	AVRG	0.167	5.4
Carbon Disulfide	AVRG	1.587	3.6
Trans-1,2-Dichloroethene	AVRG	0.532	3.3
Vinyl Acetate	AVRG	1.058	7.8
1,1-Dichloroethane	AVRG	0.962	3.3
2-Butanone	AVRG	0.253	4.7
2,2-Dichloropropane	AVRG	0.820	5.9
Cis-1,2-Dichloroethene	AVRG	0.563	2.6
Chloroform	AVRG	0.943	3.2
Bromochloromethane	AVRG	0.259	3.0
1,1,1-Trichloroethane	AVRG	0.882	2.4
1,1-Dichloropropene	AVRG	0.509	2.4
Carbon Tetrachloride	AVRG	0.468	6.0
1,2-Dichloroethane	AVRG	0.490	3.4
Benzene	AVRG	1.426	3.2
Trichloroethene	AVRG	0.366	5.0
1,2-Dichloropropane	AVRG	0.360	3.9
Bromodichloromethane	AVRG	0.459	6.9
Dibromomethane	AVRG	0.197	4.1
2-Chloroethyl Vinyl Ether	AVRG	0.140	19.1
4-Methyl-2-Pentanone	AVRG	0.136	7.7
Cis 1,3-dichloropropene	AVRG	0.566	7.2
Toluene	AVRG	0.918	3.2
Trans 1,3-Dichloropropene	AVRG	0.523	9.2
2-Hexanone	AVRG	0.274	6.4

<- Indicates value outside QC limits:
(%RSD < 20% or R² > 0.990)

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R^2
1,1,2-Trichloroethane	AVRG	0.302	4.2
1,3-Dichloropropane	AVRG	0.576	1.9
Tetrachloroethene	AVRG	0.408	1.8
Chlorodibromomethane	AVRG	0.326	13.3
1,2-Dibromoethane	AVRG	0.306	5.6
Chlorobenzene	AVRG	1.054	4.1
Ethyl Benzene	AVRG	0.571	3.4
1,1,1,2-Tetrachloroethane	AVRG	0.353	7.4
m,p-xylene	AVRG	0.694	2.7
o-Xylene	AVRG	0.681	2.9
Styrene	AVRG	1.127	8.1
Bromoform	AVRG	0.344	19.3
1,1,2,2-Tetrachloroethane	AVRG	0.685	6.4
1,2,3-Trichloropropane	AVRG	0.233	3.4
Trans-1,4-Dichloro 2-Butene	2ORDR		0.9976
N-Propyl Benzene	AVRG	3.527	4.5
Bromobenzene	AVRG	0.777	3.3
Isopropyl Benzene	AVRG	2.571	2.9
2-Chloro Toluene	AVRG	2.481	2.5
4-Chloro Toluene	AVRG	2.307	3.1
T-Butyl Benzene	AVRG	2.354	2.5
1,3,5-Trimethyl Benzene	AVRG	2.571	2.9
1,2,4-Trimethylbenzene	AVRG	2.442	7.4
S-Butyl Benzene	AVRG	3.237	4.8
4-Isopropyl Toluene	AVRG	2.649	8.8
1,3-Dichlorobenzene	AVRG	1.562	3.0
1,4-Dichlorobenzene	AVRG	1.625	2.8
N-Butyl Benzene	AVRG	2.046	14.2
1,2-Dichlorobenzene	AVRG	1.540	3.3
1,2-Dibromo 3-Chloropropane	AVRG	0.139	12.2
1,2,4-Trichlorobenzene	AVRG	0.947	11.4
Hexachloro 1,3-Butadiene	AVRG	0.568	3.0
Naphthalene	AVRG	1.904	10.1
1,2,3-Trichlorobenzene	AVRG	0.833	10.3
Dichlorodifluoromethane	AVRG	0.474	6.8
Methyl tert butyl ether	AVRG	1.451	4.3

<- Indicates value outside QC limits:
(%RSD < 20% or R^2 > 0.990)

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R ²
d4-1,2-Dichloroethane	AVRG	0.581	0.9
d8-Toluene	AVRG	1.229	1.0
4-Bromofluorobenzene	AVRG	0.556	1.5
d4-1,2-Dichlorobenzene	AVRG	0.908	0.7
Dibromofluoromethane	AVRG	0.458	2.6

<- Indicates value outside QC limits:
(%RSD < 20% or R² > 0.990)

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Cont. Calib. Date: 09/26/11

Init. Calib. Date: 09/21/11

Cont. Calib. Time: 0939

COMPOUND	Cal Amt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
Chloromethane	0.811	0.6624	0.100	AVRG	-18.3
Vinyl Chloride	0.641	0.6043	0.010	AVRG	-5.7
Bromomethane	0.327	0.2645	0.010	AVRG	-19.1
Chloroethane	0.391	0.3501	0.010	AVRG	-10.5
Trichlorofluoromethane	0.722	0.7138	0.010	AVRG	-1.1
Acrolein	0.063	0.0569	0.010	AVRG	-9.7
112Trichloro122Trifluoroetha	0.470	0.4519	0.010	AVRG	-3.8
Acetone	0.173	0.1353	0.010	AVRG	-21.8
1,1-Dichloroethene	0.491	0.4429	0.010	AVRG	-9.8
Bromoethane	0.322	0.2964	0.010	AVRG	-8.0
Iodomethane	0.532	0.3701	0.010	AVRG	-30.4
Methylene Chloride	0.542	0.4977	0.010	AVRG	-8.2
Acrylonitrile	0.167	0.1518	0.010	AVRG	-9.1
Carbon Disulfide	1.587	1.5028	0.010	AVRG	-5.3
Trans-1,2-Dichloroethene	0.532	0.5005	0.010	AVRG	-5.9
Vinyl Acetate	1.058	0.9871	0.010	AVRG	-6.7
1,1-Dichloroethane	0.962	0.8756	0.100	AVRG	-9.0
2-Butanone	0.253	0.2338	0.010	AVRG	-7.6
2,2-Dichloropropane	0.820	0.7671	0.010	AVRG	-6.4
Cis-1,2-Dichloroethene	0.563	0.5241	0.010	AVRG	-6.9
Chloroform	0.943	0.8900	0.010	AVRG	-5.6
Bromochloromethane	0.259	0.2441	0.010	AVRG	-5.8
1,1,1-Trichloroethane	0.882	0.8295	0.010	AVRG	-6.0
1,1-Dichloropropene	0.509	0.4885	0.010	AVRG	-4.0
Carbon Tetrachloride	0.468	0.4581	0.010	AVRG	-2.1
1,2-Dichloroethane	0.490	0.4422	0.010	AVRG	-9.8
Benzene	1.426	1.3452	0.010	AVRG	-5.7
Trichloroethene	0.366	0.3516	0.010	AVRG	-3.9
1,2-Dichloropropane	0.360	0.3312	0.010	AVRG	-8.0
Bromodichloromethane	0.459	0.4406	0.010	AVRG	-4.0
Dibromomethane	0.197	0.1840	0.010	AVRG	-6.6
2-Chloroethyl Vinyl Ether	0.140	0.1201	0.010	AVRG	-14.2
4-Methyl-2-Pentanone	0.136	0.1282	0.010	AVRG	-5.7
Cis 1,3-dichloropropene	0.566	0.5365	0.010	AVRG	-5.2
Toluene	0.918	0.8681	0.010	AVRG	-5.4
Trans 1,3-Dichloropropene	0.523	0.5003	0.010	AVRG	-4.3
2-Hexanone	0.274	0.2542	0.010	AVRG	-7.2

<- Exceeds QC limit of 20% D

* RF less than minimum RF

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Cont. Calib. Date: 09/26/11

Init. Calib. Date: 09/21/11

Cont. Calib. Time: 0939

COMPOUND	Cal Amt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
1,1,2-Trichloroethane	0.302	0.2814	0.010	AVRG	-6.8
1,3-Dichloropropane	0.576	0.5226	0.010	AVRG	-9.3
Tetrachloroethene	0.408	0.4040	0.010	AVRG	-1.0
Chlorodibromomethane	0.326	0.3289	0.010	AVRG	0.9
1,2-Dibromoethane	0.306	0.2876	0.010	AVRG	-6.0
Chlorobenzene	1.054	0.9954	0.300	AVRG	-5.6
Ethyl Benzene	0.571	0.5373	0.010	AVRG	-5.9
1,1,1,2-Tetrachloroethane	0.353	0.3456	0.010	AVRG	-2.1
m,p-xylene	0.694	0.6739	0.010	AVRG	-2.9
o-Xylene	0.681	0.6435	0.010	AVRG	-5.5
Styrene	1.127	1.1296	0.010	AVRG	0.2
Bromoform	0.344	0.3439	0.100	AVRG	-0.0
1,1,2,2-Tetrachloroethane	0.685	0.6328	0.300	AVRG	-7.6
1,2,3-Trichloropropane	0.233	0.2092	0.010	AVRG	-10.2
Trans-1,4-Dichloro 2-Butene	10.000	7.018	0.010	2ORDR	-29.8
N-Propyl Benzene	3.527	3.3874	0.010	AVRG	-4.0
Bromobenzene	0.777	0.7278	0.010	AVRG	-6.3
Isopropyl Benzene	2.571	2.4754	0.010	AVRG	-3.7
2-Chloro Toluene	2.481	2.2921	0.010	AVRG	-7.6
4-Chloro Toluene	2.307	2.1471	0.010	AVRG	-6.9
T-Butyl Benzene	2.354	2.2400	0.010	AVRG	-4.8
1,3,5-Trimethyl Benzene	2.571	2.4754	0.010	AVRG	-3.7
1,2,4-Trimethylbenzene	2.442	2.4454	0.010	AVRG	0.1
S-Butyl Benzene	3.237	3.1945	0.010	AVRG	-1.3
4-Isopropyl Toluene	2.649	2.7596	0.010	AVRG	4.2
1,3-Dichlorobenzene	1.562	1.4829	0.010	AVRG	-5.1
1,4-Dichlorobenzene	1.625	1.5263	0.010	AVRG	-6.1
N-Butyl Benzene	2.046	2.1562	0.010	AVRG	5.4
1,2-Dichlorobenzene	1.540	1.4386	0.010	AVRG	-6.6
1,2-Dibromo 3-Chloropropane	0.139	0.1218	0.010	AVRG	-12.4
1,2,4-Trichlorobenzene	0.947	0.9954	0.010	AVRG	5.1
Hexachloro 1,3-Butadiene	0.568	0.5674	0.010	AVRG	-0.1
Naphthalene	1.904	2.0082	0.010	AVRG	5.5
1,2,3-Trichlorobenzene	0.833	0.8534	0.010	AVRG	2.4
Dichlorodifluoromethane	0.474	0.4563	0.010	AVRG	-3.7
Methyl tert butyl ether	1.451	1.3215	0.010	AVRG	-8.9

<- Exceeds QC limit of 20% D
* RF less than minimum RF

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Cont. Calib. Date: 09/26/11

Init. Calib. Date: 09/21/11

Cont. Calib. Time: 0939

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
d4-1,2-Dichloroethane	0.581	0.5513	0.010	AVRG	-5.1
d8-Toluene	1.229	1.2170	0.010	AVRG	-1.0
4-Bromofluorobenzene	0.556	0.5536	0.010	AVRG	-0.4
d4-1,2-Dichlorobenzene	0.908	0.9103	0.010	AVRG	0.2
Dibromofluoromethane	0.458	0.4536	0.010	AVRG	-1.0

<- Exceeds QC limit of 20% D

* RF less than minimum RF

8A
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIGA

Ical Midpoint ID: 1000727

Ical Date: 07/27/11

Instrument ID: NT5

Project Run Date: 09/21/11

	IS1 (PFB) AREA #	RT #	IS2 (DFB) AREA #	RT #	IS3 (CLB) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	319331	4.77	421917	5.21	402459	7.67
UPPER LIMIT	638662	5.27	843834	5.71	804918	8.17
LOWER LIMIT	159666	4.27	210958	4.71	201230	7.17
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 LCS0921	333608	4.77	454880	5.21	420554	7.67
02 LCS0921	316220	4.77	427131	5.21	395480	7.67
03 MB0921	297550	4.77	419779	5.21	387163	7.66
04 FRP-091911-0	243958	4.77	318525	5.21	289593	7.66
05 FRP-091911-0	259731	4.77	365931	5.21	382933	7.66
06 FRP-091911-0	292481	4.85	462562	5.28	394770	7.75
07 FRP-091911-0	293301	4.77	385524	5.21	364245	7.66
08 FRP-091911-0	265354	4.77	356334	5.21	338365	7.66
09 TRIP BLANKS	258822	4.77	347379	5.21	328343	7.66
10 FRP-092011-0	264434	4.77	356649	5.21	320607	7.67
11 FRP-092011-0	258738	4.77	345053	5.21	317057	7.67
12 TRIP BLANKS	255052	4.77	337989	5.21	315968	7.66
13						
14						
15						
16						
17						
18						
19						
20						
21						
22						

IS1 (PFB) = Pentafluorobenzene

IS2 (DFB) = 1,4-Difluorobenzene

IS3 (CLB) = d5-Chlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint

AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint

RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint

RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

* Values outside of QC limits.

8A
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIGA

Ical Midpoint ID: 1000727

Ical Date: 07/27/11

Instrument ID: NT5

Project Run Date: 09/21/11

	IS4 (DCB) AREA #	RT #	AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	245371	9.72				
UPPER LIMIT	490742	10.22				
LOWER LIMIT	122686	9.22				
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 LCS0921	251539	9.72				
02 LCS0921	223163	9.72				
03 MB0921	200860	9.72				
04 FRP-091911-0	152885	9.72				
05 FRP-091911-0	206198	9.72				
06 FRP-091911-0	152847	9.76				
07 FRP-091911-0	214077	9.72				
08 FRP-091911-0	186056	9.72				
09 TRIP BLANKS	174317	9.72				
10 FRP-092011-0	179204	9.72				
11 FRP-092011-0	173682	9.72				
12 TRIP BLANKS	167717	9.72				
13						
14						
15						
16						
17						
18						
19						
20						
21						
22						

IS4 (DCB) = d4-1,4-Dichlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

* Values outside of QC limits.

8A
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIGA

Ical Midpoint ID: 10_00921

Ical Date: 09/21/11

Instrument ID: NT2

Project Run Date: 09/26/11

	IS1 (PFB) AREA #	RT #	IS2 (DFB) AREA #	RT #	IS3 (CLB) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	331438	5.46	505777	5.86	491374	7.92
UPPER LIMIT	662876	5.96	1011554	6.36	982748	8.42
LOWER LIMIT	165719	4.96	252888	5.36	245687	7.42
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 LCS0926	302141	5.46	463384	5.86	458683	7.92
02 LCS0926	294107	5.46	447017	5.86	452687	7.92
03 MB0926	310738	5.46	468089	5.86	469380	7.92
04 TRIP BLANKS	306553	5.46	459320	5.86	459420	7.92
05 FRP-091911-0	321519	5.46	480140	5.86	472582	7.92
06 FRP-091911-0	312530	5.46	469646	5.85	462564	7.92
07 FRP-091911-0	318130	5.46	478761	5.85	472117	7.92
08 FRP-091911-0	312866	5.46	470323	5.85	465816	7.92
09 FRP-092011-0	305691	5.46	455665	5.86	454285	7.92
10 FRP-092011-0	297479	5.46	445846	5.86	448317	7.92
11 FRP-092011-0	304173	5.46	455255	5.86	455236	7.92
12 FRP-092111-0	317417	5.46	475118	5.85	466598	7.92
13 FRP-092111-0	321297	5.46	479180	5.85	472408	7.92
14 FRP-092111-0	318994	5.46	480834	5.85	469531	7.92
15 FRP-092111-0	324022	5.46	490358	5.85	479960	7.92
16 FRP-092111-0	319766	5.46	483678	5.85	474939	7.92
17 FRP-092111-0	299278	5.46	446469	5.86	445768	7.92
18 FRP-091911-0	315458	5.46	478667	5.85	464572	7.92
19 FRP-091911-0	324929	5.46	492284	5.85	477416	7.92
20						
21						
22						

IS1 (PFB) = Pentafluorobenzene

IS2 (DFB) = 1,4-Difluorobenzene

IS3 (CLB) = d5-Chlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint

AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint

RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint

RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

* Values outside of QC limits.

8A
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIGA

Ical Midpoint ID: 10_00921

Ical Date: 09/21/11

Instrument ID: NT2

Project Run Date: 09/26/11

	IS4 (DCB) AREA #	RT #	AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	306350	9.62				
UPPER LIMIT	612700	10.12				
LOWER LIMIT	153175	9.12				
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 LCS0926	286825	9.61				
02 LCS0926	286948	9.61				
03 MB0926	284610	9.61				
04 TRIP BLANKS	282436	9.61				
05 FRP-091911-0	286008	9.61				
06 FRP-091911-0	277671	9.61				
07 FRP-091911-0	278594	9.61				
08 FRP-091911-0	277588	9.61				
09 FRP-092011-0	266535	9.61				
10 FRP-092011-0	273338	9.61				
11 FRP-092011-0	277228	9.61				
12 FRP-092111-0	278502	9.61				
13 FRP-092111-0	275481	9.61				
14 FRP-092111-0	270111	9.61				
15 FRP-092111-0	277630	9.61				
16 FRP-092111-0	279134	9.61				
17 FRP-092111-0	275086	9.61				
18 FRP-091911-0	284602	9.61				
19 FRP-091911-0	279507	9.61				
20						
21						
22						

IS4 (DCB) = d4-1,4-Dichlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

* Values outside of QC limits.

**Metals Analysis
Report and Summary QC Forms**

ARI Job ID: TN42, TN43

Cover Page**INORGANIC ANALYSIS DATA PACKAGE**

ANALYTICAL
RESOURCES
INCORPORATED

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

CLIENT ID	ARI ID	ARI LIMS ID	REPREP
FRP-092111-001	TN42A	11-20663	
FRP-092111-001D	TN42ADUP	11-20663	
FRP-092111-001S	TN42ASPK	11-20663	
FRP-092111-002	TN42B	11-20664	
PBW	TN42MB1	11-20664	
LCSW	TN42MB1SPK	11-20664	
FRP-092111-003	TN42C	11-20665	
FRP-092111-004	TN42D	11-20666	
FRP-092111-005	TN42E	11-20667	
FRP-092111-006	TN42F	11-20668	

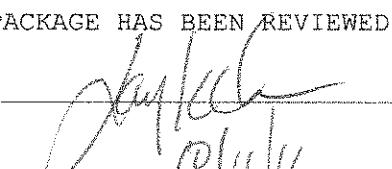
Were ICP interelement corrections applied ? Yes/No YES

Were ICP background corrections applied ? Yes/No YES

If yes - were raw data generated before application of background corrections ? Yes/No NO

Comments: _____

THIS DATA PACKAGE HAS BEEN REVIEWED AND AUTHORIZED FOR RELEASE BY:

Signature:  Name: Jay KuhnDate: 10/11/11 Title: Inorganics Director

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092111-001
SAMPLE

Lab Sample ID: TN42A

LIMS ID: 11-20663

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/21/11

Date Received: 09/21/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/23/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	48.4	
200.8	09/23/11	200.8	10/06/11	7440-38-2	Arsenic	0.00024	0.001	0.019	
3010A	09/23/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/23/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.182	
3010A	09/23/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.154	
200.8	09/23/11	200.8	10/06/11	7439-92-1	Lead	0.000230	0.0005	0.0108	
3010A	09/23/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.07	
3010A	09/23/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.391	
3010A	09/23/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.20	

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092111-002

SAMPLE

Lab Sample ID: TN42B

LIMS ID: 11-20664

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/21/11

Date Received: 09/21/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/23/11	6010B	09/27/11	7429-90-5	Aluminum	0.051	0.1	321	
200.8	09/23/11	200.8	10/06/11	7440-38-2	Arsenic	0.000120	0.0005	0.107	
3010A	09/23/11	6010B	09/27/11	7440-43-9	Cadmium	0.00036	0.004	0.006	
3010A	09/23/11	6010B	09/27/11	7440-47-3	Chromium	0.0025	0.01	0.53	
3010A	09/23/11	6010B	09/27/11	7440-50-8	Copper	0.00184	0.004	1.88	
200.8	09/23/11	200.8	10/06/11	7439-92-1	Lead	0.000115	0.0002	0.361	
3010A	09/23/11	6010B	09/27/11	7440-02-0	Nickel	0.0077	0.02	0.34	
3010A	09/23/11	6010B	09/27/11	7782-49-2	Selenium	0.010	0.1	0.1	U
3010A	09/23/11	6010B	09/27/11	7440-28-0	Thallium	0.006	0.1	0.1	U
3010A	09/23/11	6010B	09/27/11	7440-62-2	Vanadium	0.00054	0.006	1.32	
3010A	09/23/11	6010B	09/27/11	7440-66-6	Zinc	0.0029	0.02	1.27	

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

INORGANICS ANALYSIS DATA SHEET
TOTAL METALS

Page 1 of 1

Sample ID: FRP-092111-003
SAMPLE

Lab Sample ID: TN42C

LIMS ID: 11-20665

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/21/11

Date Received: 09/21/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/23/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	17.1	
200.8	09/23/11	200.8	10/06/11	7440-38-2	Arsenic	0.000120	0.0005	0.0085	
3010A	09/23/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/23/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.082	
3010A	09/23/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.137	
200.8	09/23/11	200.8	10/06/11	7439-92-1	Lead	0.000115	0.0002	0.0152	
3010A	09/23/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.02	
3010A	09/23/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.398	
3010A	09/23/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.17	

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092111-004

SAMPLE

Lab Sample ID: TN42D

LIMS ID: 11-20666

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/21/11

Date Received: 09/21/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/23/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	17.0	
200.8	09/23/11	200.8	10/06/11	7440-38-2	Arsenic	0.000120	0.0005	0.0090	
3010A	09/23/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/23/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.079	
3010A	09/23/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.128	
200.8	09/23/11	200.8	10/06/11	7439-92-1	Lead	0.000115	0.0002	0.0150	
3010A	09/23/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.02	
3010A	09/23/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U J
3010A	09/23/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.385	
3010A	09/23/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.16	

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092111-005

SAMPLE

Lab Sample ID: TN42E

LIMS ID: 11-20667

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/21/11

Date Received: 09/21/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/23/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	13.4	
200.8	09/23/11	200.8	10/06/11	7440-38-2	Arsenic	0.000120	0.0005	0.0103	
3010A	09/23/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/23/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.129	
3010A	09/23/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.082	
200.8	09/23/11	200.8	10/06/11	7439-92-1	Lead	0.000115	0.0002	0.0083	
3010A	09/23/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.02	
3010A	09/23/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.720	
3010A	09/23/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.09	

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

INORGANICS ANALYSIS DATA SHEET
TOTAL METALS

Page 1 of 1

Sample ID: FRP-092111-006
SAMPLE

Lab Sample ID: TN42F

LIMS ID: 11-20668

Matrix: Water

Data Release Authorized

Reported: 10/11/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/21/11

Date Received: 09/21/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/23/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	0.05	U
200.8	09/23/11	200.8	10/07/11	7440-38-2	Arsenic	0.000048	0.0002	0.0002	U
3010A	09/23/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/23/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.005	U
3010A	09/23/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.002	U
200.8	09/23/11	200.8	10/07/11	7439-92-1	Lead	0.000046	0.0001	0.0001	U
3010A	09/23/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.01	U
3010A	09/23/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.003	U
3010A	09/23/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.01	U

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

INORGANICS ANALYSIS DATA SHEET
TOTAL METALS

Page 1 of 1

Sample ID: FRP-092111-001

MATRIX SPIKE

Lab Sample ID: TN42A

LIMS ID: 11-20663

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/21/11

Date Received: 09/21/11

MATRIX SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Spike	Spike Added	% Recovery	Q
Aluminum	6010B	48.4	56.8	2.00	420%	H
Arsenic	200.8	0.019	0.046	0.025	108%	
Cadmium	6010B	0.002 U	0.539	0.500	108%	
Chromium	6010B	0.182	0.684	0.500	100%	
Copper	6010B	0.154	0.700	0.500	109%	
Lead	200.8	0.0108	0.0342	0.0250	93.6%	
Nickel	6010B	0.068	0.544	0.500	95.2%	
Selenium	6010B	0.05 U	1.42	2.00	71.0%	N
Thallium	6010B	0.05 U	2.05	2.00	102%	
Vanadium	6010B	0.391	0.857	0.500	93.2%	
Zinc	6010B	0.197	0.697	0.500	100%	

Reported in mg/L

N-Control Limit Not Met

H-% Recovery Not Applicable, Sample Concentration Too High

NA-Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%

INORGANICS ANALYSIS DATA SHEET
TOTAL METALS

Page 1 of 1

 Sample ID: FRP-092111-001
 DUPLICATE

Lab Sample ID: TN42A

LIMS ID: 11-20663

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/21/11

Date Received: 09/21/11

MATRIX DUPLICATE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Aluminum	6010B	48.4	49.9	3.1%	+/- 20%	
Arsenic	200.8	0.019	0.018	5.4%	+/- 20%	
Cadmium	6010B	0.002 U	0.002 U	0.0%	+/- 0.002	L
Chromium	6010B	0.182	0.181	0.6%	+/- 20%	
Copper	6010B	0.154	0.153	0.7%	+/- 20%	
Lead	200.8	0.0108	0.0108	0.0%	+/- 20%	
Nickel	6010B	0.07	0.07	0.0%	+/- 20%	
Selenium	6010B	0.05 U	0.05 U	0.0%	+/- 0.05	L
Thallium	6010B	0.05 U	0.05 U	0.0%	+/- 0.05	L
Vanadium	6010B	0.391	0.383	2.1%	+/- 20%	
Zinc	6010B	0.20	0.20	0.0%	+/- 20%	

Reported in mg/L

*-Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: TN42LCS

LIMS ID: 11-20664

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: NA

Date Received: NA

BLANK SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Spike Found	Spike Added	% Recovery	Q
Aluminum	6010B	2.05	2.00	102%	
Arsenic	200.8	0.0257	0.0250	103%	
Cadmium	6010B	0.514	0.500	103%	
Chromium	6010B	0.502	0.500	100%	
Copper	6010B	0.506	0.500	101%	
Lead	200.8	0.0254	0.0250	102%	
Nickel	6010B	0.48	0.50	96.0%	
Selenium	6010B	1.86	2.00	93.0%	
Thallium	6010B	1.87	2.00	93.5%	
Vanadium	6010B	0.494	0.500	98.8%	
Zinc	6010B	0.48	0.50	96.0%	

Reported in mg/L

N-Control limit not met

Control Limits: 80-120%

INORGANICS ANALYSIS DATA SHEET
TOTAL METALS

Page 1 of 1

Sample ID: METHOD BLANK

Lab Sample ID: TN42MB

LIMS ID: 11-20664

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN42-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: NA

Date Received: NA

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/23/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	0.05	U
200.8	09/23/11	200.8	10/06/11	7440-38-2	Arsenic	0.000048	0.0002	0.0002	U
3010A	09/23/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/23/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.005	U
3010A	09/23/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.002	U
200.8	09/23/11	200.8	10/06/11	7439-92-1	Lead	0.000046	0.0001	0.0001	U
3010A	09/23/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.01	U
3010A	09/23/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/23/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.003	U
3010A	09/23/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.01	U

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

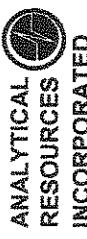
Calibration Verification

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

UNITS: ug/L



ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTv	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Aluminum	AL	ICP IP092771	2000.0	1988.30	99.4	2000.0	2017.69	100.9	2040.49	102.0	1970.63	98.5	1963.83	98.2	1971.51	98.6	
Arsenic	AS	PMS MS100682	50.0	50.04	100.1	50.0	49.47	98.9	49.11	98.2	49.56	99.1	50.17	100.3	49.55	99.1	
Cadmium	CD	ICP IP092771	1000.0	1009.47	100.9	1000.0	1016.89	101.7	1035.92	103.6	1021.69	102.2	1002.41	100.2	1002.51	100.3	
Chromium	CR	ICP IP092771	1000.0	992.76	99.3	1000.0	997.19	99.7	1017.50	101.8	984.05	98.4	977.84	97.8	984.62	98.5	
Copper	CU	ICP IP092771	1000.0	970.43	97.0	1000.0	980.82	98.1	1000.32	100.0	990.53	99.1	973.04	97.3	966.19	96.6	
Lead	PB	PMS MS100682	50.0	49.54	99.1	50.0	49.07	98.1	48.89	97.8	48.84	97.7	48.99	98.0	49.54	99.1	
Nickel	NI	ICP IP092771	1000.0	987.11	98.7	1000.0	977.68	97.8	992.25	99.2	949.35	94.9	945.56	94.6	950.03	95.0	
Selenium	SE	ICP IP092771	2000.0	1889.99	94.5	2000.0	1882.97	94.1	1895.26	94.8	1850.95	92.5	1809.01	90.5	1803.80	90.2	
Thallium	Tl	ICP IP092771	2000.0	1888.64	94.4	2000.0	1880.70	94.0	1897.32	94.9	1840.73	92.0	1800.53	90.0	1795.71	89.8	
Vanadium	V	ICP IP092771	1000.0	936.07	93.6	1000.0	940.86	94.1	959.18	95.9	945.69	94.6	930.59	93.1	927.77	92.8	
Zinc	ZN	ICP IP092771	1000.0	978.61	97.9	1000.0	975.28	97.5	996.03	99.6	962.06	96.2	958.41	95.8	969.58	97.0	

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

742-66988

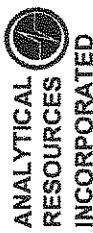
Calibration Verification

CLIENT: AMEC Geomatix

PROJECT: ERP 2011 Shoreline I

SDG: TN42

UNITS: ug/L



ANALYTE	EL	M	RUN	CCVTV	CCV6 %R	CCV7 %R	CCV8 %R	CCV9 %R	CCV10 %R	CCV11 %R
Aluminum	AL	ICP	IP092771	2000.0	1953.33	97.7	1985.46	99.3	1980.30	99.0
Arsenic	AS	PMS	MS100682	50.0	50.12	100.2	50.83	101.7	51.59	103.2
Cadmium	CD	ICP	IP092771	1000.0	1011.85	101.2	1024.67	102.5	1017.58	101.8
Chromium	CR	ICP	IP092771	1000.0	993.29	99.3	1010.51	101.1	1003.72	100.4
Copper	CU	ICP	IP092771	1000.0	964.69	96.5	977.08	97.7	971.83	97.2
Lead	PB	PMS	MS100682	50.0	48.91	97.8	49.44	98.9	49.91	99.8
Nickel	NI	ICP	IP092771	1000.0	1012.21	101.2	1032.71	103.3	1028.05	102.8
Selenium	SE	ICP	IP092771	2000.0	2030.04	101.5	2043.80	102.2	2016.20	100.8
Thallium	TL	ICP	IP092771	2000.0	2000.47	100.0	2016.53	100.8	1994.11	99.7
Vanadium	V	ICP	IP092771	1000.0	928.54	92.9	946.09	94.6	939.76	94.0
Zinc	ZN	ICP	IP092771	1000.0	994.84	99.5	1013.46	101.3	1006.94	100.7

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

Calibration Verification

CLIENT: ANEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

UNITS: ug/L



ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTv	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Arsenic	AS	PMS	MS100781	50.0	49.12	98.2	50.0	49.66	99.3	50.34	100.7	50.86	101.7				
Lead	PB	PMS	MS100781	50.0	48.89	97.8	50.0	48.79	97.6	48.49	97.0	47.95	95.9				

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

CRDL Standard

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

UNITS: ug/L

ANALYTE	EL	M	RUN	CRA/I	TV	CR-1	%R	CR-2	%R	CR-3	%R	CR-4	%R	CR-5	%R	CR-6	%R
Aluminum	AL	ICP	IP092771	50.0		53.37	106.7	49.58	99.2								
Arsenic	AS	PMS	MS100682	0.2		0.19	95.0										
Cadmium	CD	ICP	IP092771	2.0		2.27	113.5	2.13	106.5								
Chromium	CR	ICP	IP092771	5.0		4.96	99.2	4.57	91.4								
Copper	CU	ICP	IP092771	2.0		2.84	142.0	1.90	95.0								
Lead	PB	PMS	MS100682	0.1		0.11	110.0										
Nickel	NI	ICP	IP092771	10.0		9.00	90.0										
Selenium	SE	ICP	IP092771	50.0		50.50	101.0										
Thallium	TL	ICP	IP092771	50.0		47.14	94.3	48.83	97.7								
Vanadium	V	ICP	IP092771	3.0		3.10	103.3	3.01	100.3								
Zinc	ZN	ICP	IP092771	10.0		9.91	99.1	9.42	94.2								
Arsenic	AS	PMS	MS100781	0.2		0.17	85.0										
Lead	PB	PMS	MS100781	0.1		0.11	110.0										

Control Limits: no control limits have been established by the EPA at this time.

Calibration Blanks

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

UNITS: ug/L

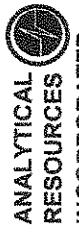
ANALYTE	EL	METH	RUN	CRDL	IDL	ICB	C	CCB1	C	CCB2	C	CCB3	C	CCB4	C	CCB5	C
Aluminum	AL	ICP	IPO92771	200.0	50.0	50.0	U										
Arsenic	AS	PMS	MS100682	10.0	0.2	0.2	U										
Cadmium	CD	ICP	IPO92771	5.0	2.0	2.0	U										
Chromium	CR	ICP	IPO92771	10.0	5.0	5.0	U										
Copper	CU	ICP	IPO92771	25.0	2.0	2.0	U										
Lead	PB	PMS	MS100682	3.0	0.1	0.1	U										
Nickel	NI	ICP	IPO92771	40.0	10.0	10.0	U										
Selenium	SE	ICP	IPO92771	5.0	50.0	50.0	U										
Thallium	TL	ICP	IPO92771	10.0	50.0	50.0	U										
Vanadium	V	ICP	IPO92771	50.0	3.0	3.0	U										
Zinc	ZN	ICP	IPO92771	20.0	10.0	10.0	U										

Calibration Blanks

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN42



UNITS: L/bn

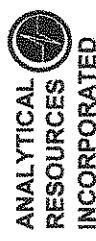
TAKE 2 : 000091

Calibration Blanks

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN42



UNITS: ug/L

ANALYTE	EL	METH	RUN	CRDL	IDL	ICB	C	CCB1	C	CCB2	C	CCB3	C	CCB4	C	CCB5	C
Arsenic	AS	PMS	MS100781	10.0	0.2	0.2	v	0.2	v	0.2	v	0.2	v	0.2	v	0.2	v
Lead	PB	PMS	MS100781	3.0	0.1	0.1	v	0.1	v	0.1	v	0.1	v	0.1	v	0.1	v

ICP Interference Check Sample

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN42



ICS SOURCE: I.V.

RUNID: IP092771

INSTRUMENT ID: OPTIMA ICP 2

UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSAI	ICSAI	%R	ICSA2	ICSA2	ICSA2	%R	ICSA3	ICSA3	%R
Aluminum	200000	200000	206000.6	201923.8	101.0	195055.3	196102.2	98.1				
Antimony	1000	37.6	979.0	97.9	26.9	950.8	95.1					
Arsenic	1000	4.3	960.0	96.0	2.6	1015.2	101.5					
Barium	1000	0.8	980.4	98.0	0.5	983.3	98.3					
Beryllium	1000	0.1	960.4	96.0	0.0	1000.1	100.0					
Boron		-6.4	-8.5	-8.5	-8.0	-10.4						
Cadmium	1000	1.4	1015.9	101.6	0.8	1013.8	101.4					
Calcium	100000	106593.1	103349.4	103.3	101843.4	100570.1	100.6					
Chromium	1000	-0.2	996.6	99.7	-0.1	987.1	98.7					
Cobalt	1000	2.7	995.0	99.5	2.8	984.5	98.5					
Copper	1000	-0.2	1025.2	102.5	-1.2	1009.5	101.0					
Iron	200000	201950.0	197833.9	98.9	188682.8	188999.9	94.5					
Lead	1000	-3.7	968.1	96.8	-3.8	1013.7	101.4					
Magnesium	100000	106013.5	98317.3	98.3	100062.5	95476.3	95.5					
Manganese	1000	0.4	926.8	92.7	0.2	930.9	93.1					
Molybdenum		6.9	6.1	5.6	6.6							
Nickel	1000	-2.3	932.0	93.2	-2.1	961.4	96.1					
Potassium		49.1	222.2	22.5	27.5	197.5						
Selenium	1000	-38.0	907.0	90.7	-33.6	977.4	97.7					
Silicon		-23.0	-23.4	-19.7	-18.8							
Silver	1000	-0.6	996.5	99.7	-0.6	1051.3	105.1					
Sodium		16.1	6.2	10.7	-1.3							
Strontium		10.6	10.1	10.3	10.2							
Thallium	1000	13.7	979.1	97.9	14.2	1031.3	103.1					
Tin		-7.3	-7.2	-7.6	-10.4							
Titanium		1.5	1.5	1.0	0.9							
Vanadium	1000	4.3	958.1	95.8	2.4	948.4	94.8					
Zinc	1000	-1.2	930.8	93.1	-0.9	943.5	94.4					

ICP Interference Check Sample



CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

ICS SOURCE: I.V.

RUNID: MS100682

INSTRUMENT ID: PE ELAN 6000

UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSA1	ICSA1	ICSA2	ICSA2	ICSA3	ICSA3	ICSA3	%R	%R	%R
Antimony			0.1	0.1								
Cadmium	20		0.0	0.0	20.0	100.0						
Chromium	20		0.8	0.8	20.5	102.5						
Cobalt	20		0.1	0.1	19.6	98.0						
Copper	20		0.5	0.5	19.4	97.0						
Iron	20000	20000	20327.5	20246.7	101.2							
Manganese	20		0.7	0.7	20.0	100.0						
Molybdenum	400	400	427.1	421.6	105.4							
Nickel	20		0.7	18.7	93.5							
Silver	20		0.0	19.3	96.5							
Vanadium			0.0	-0.5								
Zinc	20		1.1	19.8	99.0							

ICP Interference Check Sample

CLIENT: AMEC Geomatrix
PROJECT: FRP 2011 Shoreline I
SDG: TN42



TCS SOURCE: I.V.
RUNID: MS100781
INSTRUMENT ID: PE ELAN 6000
UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSA1	ICSA1	ICSA2	ICSA2	ICSA3	ICSA3	%R	%R
Antimony			0.1	0.1						
Cadmium	2.0		0.0	0.0	20.0	100.0				
Chromium	20		0.8	0.8	20.2	101.0				
Cobalt	20		0.1	0.1	18.7	93.5				
Copper	20		0.4	0.4	20.2	101.0				
Iron	20000	20000	19767.1	19767.1	19820.1	99.1				
Manganese	20		0.7	0.7	19.2	96.0				
Molybdenum	400	400	453.7	453.7	461.5	115.4				
Nickel	20		0.6	0.6	20.0	100.0				
Selenium			0.0	0.0	0.1	0.1				
Silver	20		0.0	0.0	19.3	96.5				
Vanadium			0.0	0.0	-0.4	-0.4				
Zinc	20		0.9	0.9	20.1	100.5				

**Post Digest Spike
Sample Recovery**

**ANALYTICAL
RESOURCES** 
INCORPORATED

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

ANALYSIS METHOD: ICP

SDG: TN42

UNITS: ug/L

ANALYTE	CLIENT ID	ARI ID	RUNID	SPIKED		SAMPLE RESULT C	SAMPLE RESULT C	SPIKE		MATRIX	%R
				RESULT	C			ADDED	%R		
Selenium	FRP-092111-001A	TN42APOST	IP092771	1851.06		50.00U	2000	Water	92.6		

ICP Serial Dilutions

ANALYTICAL
RESOURCES
INCORPORATED

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

ANALYSIS METHOD: ICP

SDG: TN42

UNITS: ug/L

ANALYTE	CLIENT ID	ARI ID	MATRIX	RUNID	INITIAL SAMPLE RESULT (I)	SERIAL DILUTION RESULT (S)		% DIFFER- ENCE	Q
						C	C		
Aluminum	FRP-092111-001L	TN42A-L	Water	IP092771	48407.27	49709.35		2.7	
Cadmium	FRP-092111-001L	TN42A-L	Water	IP092771	0.48 U	10.00 U			
Chromium	FRP-092111-001L	TN42A-L	Water	IP092771	181.77	185.40		2.0	
Copper	FRP-092111-001L	TN42A-L	Water	IP092771	154.47	156.05		1.0	
Nickel	FRP-092111-001L	TN42A-L	Water	IP092771	68.25	72.45 B		6.2	
Selenium	FRP-092111-001L	TN42A-L	Water	IP092771	1.97 U	250.00 U			
Thallium	FRP-092111-001L	TN42A-L	Water	IP092771	4.06 U	250.00 U			
Vanadium	FRP-092111-001L	TN42A-L	Water	IP092771	390.68	400.50		2.5	
Zinc	FRP-092111-001L	TN42A-L	Water	IP092771	196.74	203.15		3.3	

ICP Serial Dilutions

ANALYTICAL
RESOURCES
INCORPORATED

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

ANALYSIS METHOD: PMS

SDG: TN42

UNITS: ug/L

ANALYTE	CLIENT ID	ARI ID	MATRIX	RUNID	(I)	INITIAL	SERIAL	% DIFFER-	Q
						SAMPLE RESULT	DILUTION RESULT		
						C	(S)	C	Q
Arsenic	FRP-092111-001L	TN42A-L	Water	MS100682	3.73 B	3.50 B	6.2		
Lead	FRP-092111-001L	TN42A-L	Water	MS100682	2.16 B	2.25 B	4.2		

**IDLs and ICP
Linear Ranges**



CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

UNITS: ug/L

ANALYTE	EL	METH	INSTRUMENT	WAVELENGTH (nm)	GFA BACK- GROUND	ICP		RL DATE	ICP LINEAR RANGE (ug/L)	ICP LR DATE
						CLP	CRDL			
Aluminum	AL	ICP	OPTIMA ICP 2	308.22		200	50.0	4/1/2011	250000.0	8/3/2011
Arsenic	AS	PMS	PE ELAN 6000 MS	0.00		10	0.2	4/1/2011		
Cadmium	CD	ICP	OPTIMA ICP 2	228.80		5	2.0	4/1/2011	20000.0	8/3/2011
Chromium	CR	ICP	OPTIMA ICP 2	267.72		10	5.0	4/1/2011	100000.0	8/3/2011
Copper	CU	ICP	OPTIMA ICP 2	324.75		25	2.0	4/1/2011	40000.0	8/3/2011
Lead	PB	PMS	PE ELAN 6000 MS	0.00		3	0.1	4/1/2011		
Nickel	NI	ICP	OPTIMA ICP 2	231.60		40	10.0	4/1/2011	100000.0	8/3/2011
Selenium	SE	ICP	OPTIMA ICP 2	196.02		5	50.0	4/1/2011	20000.0	8/3/2011
Thallium	TL	ICP	OPTIMA ICP 2	190.86		10	50.0	4/1/2011	30000.0	8/3/2011
Vanadium	V	ICP	OPTIMA ICP 2	292.40		50	3.0	4/1/2011	50000.0	8/3/2011
Zinc	ZN	ICP	OPTIMA ICP 2	213.86		20	10.0	4/1/2011	100000.0	8/3/2011

ICP Interelement Correction Factors



CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

IEC DATE: 9/26/2011
INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELLENGTH	AL	AS	BA	BE	CA	CD	CO	CR	CU	FE
Aluminum	308.22	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Antimony	206.84	0.3618900	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	10.6149000	0.0000000	0.0000000
Arsenic	188.98	0.0473543	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.8345790	1.1215160	0.0000000	0.0000000
Barium	233.53	0.0000000	0.0000000	0.0000000	0.0000000	-0.0061516	0.0000000	-0.1892080	0.0000000	0.0000000	0.0499666
Beryllium	313.04	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Cadmium	228.80	0.0000000	5.2418600	0.0000000	0.0000000	0.0000000	0.0000000	0.1195910	0.0000000	0.0000000	0.0000000
Calcium	317.93	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.5252460	0.0000000	0.0000000
Chromium	267.72	0.0000000	0.0000000	0.0000000	0.0000000	0.020343	0.0000000	0.0000000	0.0000000	0.0000000	-0.0461558
Cobalt	228.62	0.0000000	0.0000000	0.0000000	0.1238430	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Copper	324.75	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.2279050	-0.03188969	0.0000000	-0.0514304
Iron	273.96	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-1.6163960	0.0000000	0.0000000	0.0000000
Lead	220.35	-0.2408020	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-2.3072160	1.2452600	0.0597200
Magnesium	279.08	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-1.6380600	-1.2519300	0.0000000	0.5288270
Manganese	257.61	0.0065976	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.0045330
Molybdenum	202.03	0.0000000	0.0000000	0.0000000	0.0000000	0.0103467	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Nickel	231.60	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Potassium	766.49	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Selenium	196.03	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.3514040	0.0000000	0.0000000	0.0000000
Silicon	288.16	0.0000000	0.0000000	0.0000000	0.0000000	-3.4885200	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Silver	328.07	0.0000000	0.0000000	0.0000000	0.0000000	0.0118175	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Sodium	589.59	0.0000000	0.0000000	0.0000000	0.0000000	8.8385500	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Thallium	190.80	0.1350590	0.0000000	0.0000000	0.0000000	0.0000000	1.7836600	0.3510820	0.0000000	-0.1171730	
Tin	189.93	0.0000000	0.0000000	0.0000000	0.0000000	-0.0244672	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Titanium	334.90	0.0000000	0.0000000	0.0000000	0.0000000	0.0593584	0.0000000	0.0000000	0.1632010	0.0000000	0.0000000
Vanadium	292.40	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-3.9361800	0.0000000	0.0934302	
Zinc	206.20	0.0000000	0.0000000	0.0000000	0.0135250	0.0000000	-0.1475900	0.0000000	0.0000000	0.0000000	

ICP Interelement Correction Factors

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN42



TEC DATE:

9/26/2011

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	MG	MN	MO	NI	PB	SB	TI	TL	V	ZN
Aluminum	308.22	0.0000000	0.0000000	15.3131000	0.0000000	0.0000000	1.5167500	0.0000000	17.6996000	0.0000000	0.0000000
Antimony	206.84	0.0000000	0.0000000	0.0000000	-0.4730780	0.0000000	0.0000000	-0.8897510	0.0000000	-3.3546800	0.0000000
Arsenic	188.98	0.0000000	0.0000000	2.3330800	0.0000000	0.0000000	-5.4412000	0.0000000	0.0000000	0.0000000	0.0000000
Barium	233.53	0.0000000	0.0000000	0.0000000	0.0766262	0.0000000	0.0000000	0.0000000	0.0000000	0.6419380	0.0000000
Beryllium	313.04	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.2960780	0.0000000
Cadmium	228.80	0.0000000	0.0000000	0.0000000	-0.7324130	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Calcium	317.93	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Chromium	267.72	0.0464480	0.0000000	0.1395070	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.2773470	0.0000000
Cobalt	228.62	0.0000000	0.0000000	-0.1579570	0.1588330	0.0000000	0.0000000	1.8115900	0.0000000	0.0000000	0.0000000
Copper	324.75	0.0000000	0.0000000	0.2688440	0.0000000	0.0000000	0.2461180	0.0000000	0.0000000	0.0000000	0.0000000
Iron	273.96	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	8.4403600	0.0000000
Lead	220.35	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Magnesium	279.08	0.0000000	0.0000000	-4.6256200	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Manganese	257.61	0.0039764	0.0000000	0.0000000	0.0000000	-0.2175850	0.0000000	0.0000000	0.0000000	-0.0271775	0.0000000
Molybdenum	202.03	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Nickel	231.60	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.7744280	0.0000000	0.0000000	0.0000000	0.0000000
Potassium	766.49	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Selenium	196.03	0.0780238	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Silicon	288.16	0.0000000	0.0000000	-2.7358100	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Silver	328.07	0.0000000	0.2442620	0.0000000	0.0000000	0.0000000	-0.0470302	0.0000000	-0.2758080	0.0000000	0.0000000
Sodium	589.59	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Thallium	190.80	0.0000000	-1.4179000	1.9562000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	1.2892100	0.0000000
Tin	189.93	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.5848020	-0.3044710	0.0000000	0.0000000	0.0000000
Titanium	334.90	0.0000000	0.0000000	0.9873960	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Vanadium	292.40	0.0000000	-0.1398510	-0.6804250	0.0000000	0.0000000	0.6004670	0.0000000	0.0000000	0.0000000	0.0000000
Zinc	206.20	0.0000000	0.0000000	0.2377960	0.0000000	-0.0708227	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

Preparation Log

ANALYTICAL
RESOURCES
INCORPORATED

CLIENT: AMEC Geomatrix

ANALYSIS METHOD: ICP

PROJECT: FRP 2011 Shoreline I

ARI PREP CODE: TWC

SDG: TN42

PREPDATE: 9/23/2011

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
FRP-092111-001	TN42A	0.000	50.0	50.0
FRP-092111-001D	TN42ADUP	0.000	50.0	50.0
FRP-092111-001S	TN42ASPK	0.000	50.0	50.0
FRP-092111-002	TN42B	0.000	50.0	50.0
FRP-092111-003	TN42C	0.000	50.0	50.0
FRP-092111-004	TN42D	0.000	50.0	50.0
FRP-092111-005	TN42E	0.000	50.0	50.0
FRP-092111-006	TN42F	0.000	50.0	50.0
PBW	TN42MB1	0.000	50.0	50.0
LCSW	TN42MB1SPK	0.000	50.0	50.0

Preparation Log

ANALYTICAL
RESOURCES
INCORPORATED

CLIENT: AMEC Geomatrix

ANALYSIS METHOD: PMS

PROJECT: FRP 2011 Shoreline I

ARI PREP CODE: REN

SDG: TN42

PREPDATE: 9/23/2011

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
FRP-092111-001	TN42A	0.000	50.0	25.0
FRP-092111-001D	TN42ADUP	0.000	50.0	25.0
FRP-092111-001S	TN42ASPK	0.000	50.0	25.0
FRP-092111-002	TN42B	0.000	50.0	25.0
FRP-092111-003	TN42C	0.000	50.0	25.0
FRP-092111-004	TN42D	0.000	50.0	25.0
FRP-092111-005	TN42E	0.000	50.0	25.0
FRP-092111-006	TN42F	0.000	50.0	25.0
PBW	TN42MB1	0.000	50.0	25.0
LCSW	TN42MB1SPK	0.000	50.0	25.0

Analysis Run Log

CITIENT • ANEC Geomatix

PROJECT: FBP 2011 shoreline T

SDG: TN42

INSTRUMENT ID: OPTIMA ICP 2 START DATE: 9/27/2011
RUNID: I9092771 METHOD: ICP END DATE: 9/27/2011

Analysis Run Log

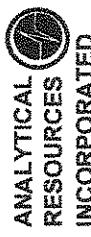
CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

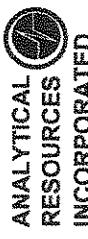
INSTRUMENT ID: OPTIMA ICP 2
RUNID: IP092771

START DATE: 9/27/2011
END DATE: 9/27/2011



CLIENT ID	ART ID	DIL.	TIME	%R	AG	AL	S	B	A	E	C	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	T1	T2	U	V	ZN
ZZZZZZ	TN19E	1.00	12495																												
LCSW	TN42MB1SPK	1.00	12540																												
CCV	CCV3	1.00	12580																												
CCB	CCB3	1.00	13014																												
S0	S0	1.00	13055																												
CCV	CCV4	1.00	13094																												
CCB	CCB4	1.00	13133																												
ZZZZZZ	TN78MB	1.00	13174																												
ZZZZZZ	TN78A	1.00	13250																												
ZZZZZZ	TN79A	1.00	13291																												
FRP-092111-003	TN42C	1.00	13332																												
FRP-092111-004	TN42D	1.00	13380																												
FRP-092111-005	TN42E	1.00	13423																												
FRP-092111-006	TN42F	1.00	13470																												
ZZZZZZ	ZZZZZZ	5.00	13511																												
FRP-092111-002	TN42B	2.00	13552																												
ZZZZZZ	TN78MBSPK	1.00	13593																												
CCV	CCV5	1.00	14033																												
CCB	CCB5	1.00	14071																												
S2	S2	1.00	14161																												
S3	S3	1.00	14173																												
CCV	CCV6	1.00	14193																												
CCB	CCB6	1.00	14232																												
FRP-092111-003	TN42C	1.00	14273																												
FRP-092111-004	TN42D	1.00	14321																												
FRP-092111-005	TN42E	1.00	14364																												
FRP-092111-006	TN42F	1.00	14411																												
ZZZZZZ	TN19APOST	5.00	14452																												
FRP-092111-002	TN42B	2.00	14493																												
CCV	CCV7	1.00	14534																												
CCB	CCB7	1.00	14573																												
CRI	CRI	1.00	15014																												
ICSA	ICSAF	1.00	15035																												
ICSA	ICSAFB	1.00	15100																												
CCV	CCV8	1.00	15135																												

Analyses Run Log



CLIENT: AMEC Geomatix

BBO Tercer Trimestre 2011 Shows Lineas T

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SDG: INDEX

INSTRUMENT ID: OBTTMA TCB 2

卷之三

KUNDI: 1E09Z/1 METHOD: ICE

START DATE: 8/27/2011

卷之三

END DATE: 9/21/2011

NAME: TRAVIS COOPER

Analysis Run Log

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

INSTRUMENT ID: PE ELAN 6000 MS
RUNID: MS100682 METHOD: PMS

START DATE: 10/6/2011
END DATE: 10/6/2011

CLIENT ID	ART ID	DIL. TIME	*R AG AL AS E BA BE CA CD CO CR CU WE HG K MG MN MO NA NT PB SB SE SI SN TI TL U V ZN
S0	S0	1.00 11460	X
ZZZZZZ	ZZZZZZ	1.00 11520	
ZZZZZZZ	ZZZZZZZ	1.00 11590	X X
S3	S3	1.00 12050	X X
S4	S4	1.00 12120	X X
S1	S1	1.00 12190	X X
S2	S2	1.00 12250	X X
ICV	MICV	1.00 12320	X X
ICB	ICB	1.00 12400	X X
CCV	MCCV1	1.00 12450	X X
CCB	CCB1	1.00 12520	X X
CRI	MCRI	1.00 12580	X X
ICSA	ICSAI	1.00 13030	X X
ICSAB	ICSABI	1.00 13100	
ZZZZZZ	LR200	1.00 13160	
ZZZZZZ	LR300	1.00 13230	X X
CCV	MCCV2	1.00 13310	
CCB	CCB2	1.00 13380	
ZZZZZZ	TM71MB	2.00 13460	
ZZZZZZ	TM67ABSPK	2.00 13520	
ZZZZZZ	TM71R	2.00 13580	
ZZZZZZ	TM71S	2.00 14050	
ZZZZZZ	TM79C	2.00 14110	
ZZZZZZ	TM67C	2.00 14180	
ZZZZZZ	TM67A	5.00 14240	
ZZZZZZ	TM67B	1000.00 14310	
ZZZZZZ	TM79H	20.00 14370	X X
ZZZZZZ	TM79J	5.00 14430	
ZZZZZZ	ZZZZZZ	1.00 14500	
CCV	MCCV3	1.00 14560	
CCB	CCB3	1.00 15030	
ZZZZZZ	TM79FDUP	1000.00 15150	
ZZZZZZ	TM79F	1000.00 15210	
ZZZZZZ	TM79FSK	1000.00 15270	
ZZZZZZ	TM79K	1000.00 15330	

Analysis Run Log

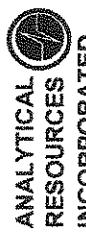
CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

INSTRUMENT ID: PE ELAN 6000 MS
RUNID: MS100682 METHOD: PMS

START DATE: 10/6/2011
END DATE: 10/6/2011



CLIENT ID	ARI ID	DLY. TIME	SR	AG AL AS B	BA BE CA CD CO CR CU FE HG K MG MN MO NA NI PB SB SE SI SN TI TL U V ZN	
ZZZZZZ	TM79G	1000.00 15390				
ZZZZZZ	TM79G	2.00 15460				
ZZZZZZ	TN00A-L	100.00 15520				
ZZZZZZ	TN00A	20.00 15590				
ZZZZZZ	TN00ADUP	20.00 16050				
ZZZZZZ	TN00ASPK	20.00 16120				
CCV	MCCV4	1.00 16220				
CCB	CCB4	1.00 16300				
ZZZZZZ	TN00MB1	2.00 16350				
ZZZZZZ	TN00MB1SPK	20.00 16460				
ZZZZZZ	TN00MB2SPK	2.00 16530				
ZZZZZZ	TN00B-1	50.00 16590				
ZZZZZZ	TN00B	10.00 17060				
ZZZZZZ	TN00BDUP	10.00 17120				
ZZZZZZ	TN00BSFK	10.00 17180				
ZZZZZZ	TN00C	20.00 17250				
ZZZZZZ	TN00D	5.00 17310				
CCV	MCCV5	1.00 17370				
CCB	CCB5	1.00 17440				
ZZZZZZ	TN19MB1	2.00 17500				
ZZZZZZ	TN19MB1SPK	2.00 17550				
ZZZZZZ	TN00E	20.00 18020				
ZZZZZZ	TN00F	5.00 18080				
ZZZZZZ	ZZZZZZ	25.00 18140				
ZZZZZZ	TN19A	5.00 18210				
ZZZZZZ	TN19ADUP	5.00 18270				
ZZZZZZ	TN19ASPK	5.00 18340				
ZZZZZZ	TN19B	5.00 18400				
ZZZZZZ	TN19C	5.00 18470				
CCV	MCCV6	1.00 18520				
CCB	CCB6	1.00 18590				
PBW	TN42MB1	2.00 19030				
LCSW	TN42MB1SPK	2.00 19100				
ZZZZZZ	TN19D	5.00 19170				

FORM XIV

Analysis Run Log



CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN42

INSTRUMENT ID: PE ELAN 6000 MS
RUNID: MS100682 METHOD: PMS

START DATE: 10/6/2011

END DATE: 10/6/2011

CLIENT ID	ARI ID	DIL.	TIME	%R	AG	AL	B	BA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	TI	TL	U	V	ZN
ZZZZZZ	TN19E	5.00	19230																													
FRP-092111-001L	TN42A-L	50.00	19300																											X		
FRP-092111-001	TN42A	10.00	19360																										X			
FRP-092111-001D	TN42ADUP	10.00	19430																										X			
FRP-092111-001S	TN42ASPK	10.00	19490																									X				
FRP-092111-002	TN42B	5.00	19550																									X				
FRP-092111-003	TN42C	5.00	20020																									X				
CCV	MCCV7	1.00	20080																									X				
CCB	CCB7	1.00	20150																									X				
ZZZZZZ	TN24MB2	20.00	20210																													
FRP-092111-004	TN24MB2SPK	20.00	20270																									X				
FRP-092111-005	TN42D	5.00	20330																									X				
FRP-092111-006	TN42E	5.00	20390																									X				
ZZZZZZ	TN42F	5.00	20460																													
ZZZZZZ	TN24A	2.00	20520																													
ZZZZZZ	TN24B	2.00	20590																													
ZZZZZZ	TN24E	2.00	21050																													
ZZZZZZ	TN24C	20.00	21120																													
ZZZZZZ	TN24D	20.00	21180																													
CCV	MCCV8	1.00	21240																									X				
CCB	CCB8	1.00	21320																								X					

Analysis Run Log

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

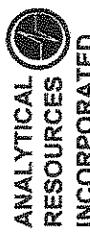
INSTRUMENT ID: PE ELAN 6000 MS

SDG: TN42

RUNID: MS100781

METHOD: PMS

START DATE: 10/7/2011
END DATE: 10/7/2011



CLIENT ID	ARI ID	DIL.	TIME	%R	AG	AL	AS	B	BA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	TI	TL	U	V	ZEN
S0	S0		1.00 08520																														
S1	S1		1.00 08580																														
S2	S2		1.00 09040																														
S3	S3		1.00 09110																														
S4	S4		1.00 09170																														
ZZZZZZ	Rinse Samp1		1.00 09240																														
ICV	MICV		1.00 09320																														
ICB	ICB		1.00 09390																														
CCV	MCCV1		1.00 09440																														
CCB	CCB1		1.00 09510																														
CRI	MCRI		1.00 09570																														
ICSA	ICSAI		1.00 10030																														
ICSAB	ICSABI		1.00 10090																														
ZZZZZZ	ZZZZZZ		1.00 10160																														
ZZZZZZ	LR200		1.00 10240																														
ZZZZZZ	LR300		1.00 10320																														
CCV	MCCV2		1.00 10390																														
CCB	CCB2		1.00 10460																														
ZZZZZZ	TN63MB2		2.00 10560																														
ZZZZZZ	TN00E		2.00 11020																														
FRP-092111-006	TN42F		2.00 11070																														
ZZZZZZ	TN19E		2.00 11130																														
ZZZZZZ	ZZZZZZ		50.00 11190																														
ZZZZZZ	TN19A		10.00 11250																														
ZZZZZZ	TN19ADUP		10.00 11310																														
ZZZZZZ	TN19ASPK		10.00 11380																														
ZZZZZZ	TN24A		100.00 11440																														
CCV	TN24A		2.00 11510																														
CCB	MCCV3		1.00 11580																														
CCB	CCB3		1.00 12050																														

Mercury Analysis
Report and Summary QC Forms

ARI Job ID: TN42, TN43

Cover Page**INORGANIC ANALYSIS DATA PACKAGE**

CLIENT: AMEC Geomatrix

PROJECT: FRP Shoreline Invest

SDG: TN43

CLIENT ID	ARI ID	ARI LIMS ID	REPREP
FRP-092111-001	TN43A	11-20670	
FRP-092111-001D	TN43ADUP	11-20670	
FRP-092111-001S	TN43ASPK	11-20670	
FRP-092111-002	TN43B	11-20671	
PBW	TN43MB1	11-20671	
LCSW	TN43MB1SPK	11-20671	
FRP-092111-003	TN43C	11-20672	
FRP-092111-004	TN43D	11-20673	
FRP-092111-005	TN43E	11-20674	
FRP-092111-006	TN43F	11-20675	
FRP-092111-001	TN43RA	11-20670	R
FRP-092111-001D	TN43RADUP	11-20670	R
FRP-092111-001S	TN43RASPK	11-20670	R
FRP-092111-002	TN43RB	11-20671	R
PBW	TN43RMB1	11-20671	R
LCSW	TN43RMB1SPK	11-20671	R
FRP-092111-003	TN43RC	11-20672	R
FRP-092111-004	TN43RD	11-20673	R
FRP-092111-005	TN43RE	11-20674	R
FRP-092111-006	TN43RF	11-20675	R

Were ICP interelement corrections applied ? Yes/No YES

Were ICP background corrections applied ? Yes/No YES

If yes - were raw data generated before application of background corrections ? Yes/No NO

Comments: _____

THIS DATA PACKAGE HAS BEEN REVIEWED AND AUTHORIZED FOR RELEASE BY:

Signature: Jay Kuhn Name: Jay KuhnDate: 10/11/11 Title: Inorganics Director

INORGANICS ANALYSIS DATA SHEET
Total Mercury by Method SW7470A

ANALYTICAL
RESOURCES
INCORPORATED

Data Release Authorized: *[Signature]*
Reported: 10/07/11
Date Received: 09/21/11
Page 1 of 1

QC Report No: TN43-AMEC Geomatrix
Project: FRP Shoreline Investigation
8769

Client/ ARI ID	Date Sampled	Matrix	Prep Date Anal Date	RL	Result
FRP-092111-001 TN43A 11-20670	09/21/11	Water	10/04/11 10/06/11	20.0	136 J
FRP-092111-002 TN43B 11-20671	09/21/11	Water	10/04/11 10/06/11	400	9,760 J
FRP-092111-003 TN43C 11-20672	09/21/11	Water	10/04/11 10/06/11	400	400 U
FRP-092111-004 TN43D 11-20673	09/21/11	Water	10/04/11 10/06/11	400	558 J
FRP-092111-005 TN43E 11-20674	09/21/11	Water	10/04/11 10/06/11	400	820 J
FRP-092111-006 TN43F 11-20675	09/21/11	Water	10/04/11 10/06/11	20.0	20.0 U
MB-100411 Method Blank	NA	Water	10/04/11 10/06/11	20.0	20.0 U

Reported in ng/L

RL-Analytical reporting limit
U-Undetected at reported detection limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092111-001

MATRIX SPIKE

Lab Sample ID: TN43A

LIMS ID: 11-20670

Matrix: Water

Data Release Authorized:

Reported: 10/07/11

QC Report No: TN43-AMEC Geomatrix

Project: FRP Shoreline Investigation

8769

Date Sampled: 09/21/11

Date Received: 09/21/11

MATRIX SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Spike	Spike Added	% Recovery	Q
Mercury	7470A	136	311	100	175%	N

Reported in ng/L

N=Control Limit Not Met

H=% Recovery Not Applicable, Sample Concentration Too High

NA=Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Lab Sample ID: TN43A

LIMS ID: 11-20670

Matrix: Water

Data Release Authorized:

Reported: 10/07/11

Sample ID: FRP-092111-001

DUPLICATE

QC Report No: TN43-AMEC Geomatrix

Project: FRP Shoreline Investigation

8769

Date Sampled: 09/21/11

Date Received: 09/21/11

MATRIX DUPLICATE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Mercury	7470A	136	307	77.2%	+/- 20%	*

Reported in ng/L

*-Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Lab Sample ID: TN43LCS

LIMS ID: 11-20671

Matrix: Water

Data Release Authorized:

Reported: 10/07/11

Sample ID: LAB CONTROL

QC Report No: TN43-AMEC Geomatrix

Project: FRP Shoreline Investigation

8769

Date Sampled: NA

Date Received: NA

BLANK SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Spike Found	Spike Added	% Recovery	Q
Mercury	7470A	201	200	100%	

Reported in ng/L

N-Control limit not met

Control Limits: 80-120%

Calibration Verification

CLIENT: AMEC Geomatix

PROJECT: FRP Shoreline Invest

SDG: TN43

UNITS: ng/L



ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTV	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Mercury	HG	CVL	HG100601	500.0	503.00	100.6	500.0	511.00	102.2	516.00	103.2	517.00	103.4	517.00	103.4	517.00	103.4

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

CRDL Standard

CLIENT: AMEC Geomatix

PROJECT: FRP Shoreline Invest

SDG: TN43

ANALYTE	EL	M	RUN	CRA/I	TV	CR-1	%R	CR-2	%R	CR-3	%R	CR-4	%R	CR-5	%R	CR-6	%R
Mercury	HG	CYL	HG100601	20.0	20.00	100.0											

UNITS: ng/L

Control Limits: no control limits have been established by the EPA at this time.



Calibration Blanks

CLIENT: AMEC Geomatix

PROJECT: FRP Shoreline Invest

SDG: TN43



UNITS: ng/L

ANALYTE	EL	METH	RUN	CRDL	IDL	ICB	C	CCB1	C	CCB2	C	CCB3	C	CCB4	C	CCB5	C
Mercury	HG	CVL	HG100601	25.0	20.0	20.0	U										

**IDLs and ICP
Linear Ranges**

**ANALYTICAL
RESOURCES
INCORPORATED**

CLIENT: AMEC Geomatrix

IDL DATE: 4/1/2011

PROJECT: FRP Shoreline Invest

ICP LR DATE:

SDG: TN43

UNITS: ng/L

ANALYTE	EL	METH	INSTRUMENT	WAVELENGTH (nm)	GFA BACK- GROUND	CLP CRDL	RL	ICP LINEAR RANGE (ng/L)
Mercury	HG	CVL	CETAC MERCURY	253.70		25	20.0	

Preparation Log

ANALYTICAL
RESOURCES
INCORPORATED

CLIENT: AMEC Geomatrix

ANALYSIS METHOD: CVL

PROJECT: FRP Shoreline Invest

ARI PREP CODE: TLM

SDG: TN43

PREPDATE: 9/23/2011

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
FRP-092111-001	TN43A	0.000	20.0	20.0
FRP-092111-001D	TN43ADUP	0.000	20.0	20.0
FRP-092111-001S	TN43ASPK	0.000	20.0	20.0
FRP-092111-002	TN43B	0.000	20.0	20.0
FRP-092111-003	TN43C	0.000	20.0	20.0
FRP-092111-004	TN43D	0.000	20.0	20.0
FRP-092111-005	TN43E	0.000	20.0	20.0
FRP-092111-006	TN43F	0.000	20.0	20.0
PBW	TN43MB1	0.000	20.0	20.0
LCSW	TN43MB1SPK	0.000	20.0	20.0

Preparation Log



CLIENT: AMEC Geomatrix

ANALYSIS METHOD: CVL

PROJECT: FRP Shoreline Invest

ARI PREP CODE: TLM

SDG: TN43

PREPDATE: 10/4/2011

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
FRP-092111-001	TN43RA	0.000	20.0	20.0
FRP-092111-001D	TN43RADUP	0.000	20.0	20.0
FRP-092111-001S	TN43RASPK	0.000	20.0	20.0
FRP-092111-002	TN43RB	0.000	1.0	20.0
FRP-092111-003	TN43RC	0.000	1.0	20.0
FRP-092111-004	TN43RD	0.000	1.0	20.0
FRP-092111-005	TN43RE	0.000	1.0	20.0
FRP-092111-006	TN43RF	0.000	20.0	20.0
PBW	TN43RMB1	0.000	20.0	20.0
LCSW	TN43RMB1SPK	0.000	20.0	20.0

Analysis Run Log

CLIENT: AMEC Geomatix

PROJECT: FRP Shoreline Invest

SDG: TN43

INSTRUMENT ID: CETAC MERCURY
RUNID: HG100601

START DATE: 10/6/2011
END DATE: 10/6/2011

CLIENT ID	ART ID	DIL.	TIME	8R	AG	AL	AS	B	EA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	T1	T1L	U	V	ZN
SO	SO		1.00	11362																													
S20	S20		1.00	11390																													
S50	S50		1.00	11414																													
S100	S100		1.00	11442																													
S200	S200		1.00	11470																													
S400	S400		1.00	11495																													
S1000	S1000		1.00	11523																													
ICV	AICV		1.00	11595																													
ICB	ICB		1.00	12023																													
CCV	ACCVI		1.00	12051																													
CCB	CCB1		1.00	12080																													
CRA	CRA		1.00	12104																													
ZZZZZZ	TN03RMH1SPK		1.00	12160																													
ZZZZZZ	TN03RA		1.00	12184																													
ZZZZZZ	TN03RADUP		1.00	12212																													
ZZZZZZ	TN03RASPK		1.00	12240																													
ZZZZZZ	TN03RB		1.00	12265																													
ZZZZZZ	TN03RC		1.00	12293																													
ZZZZZZ	TN03RD		1.00	12321																													
ZZZZZZ	TN03RE		1.00	12345																													
CCV	ACCV2		1.00	12374																													
CCB	CCB2		1.00	12402																													
PBW	TN43RMH1		1.00	12431																													
LCSW	TN43RMH1SPK		1.00	12455																													
FRP-092111-001	TN43RA		1.00	12483																													
FRP-092111-001D	TN43RADUP		1.00	12511																													
FRP-092111-001S	TN43RASPK		1.00	12535																													
FRP-092111-002	TN43RB		1.00	12563																													
FRP-092111-003	TN43RC		1.00	12591																													
FRP-092111-004	TN43RD		1.00	13015																													
FRP-092111-005	TN43RE		1.00	13044																													
FRP-092111-006	TN43RF		1.00	13072																													
CCV	ACCV3		1.00	13100																													
CCB	CCB3		1.00	13125																													

Analysis Run Log



CLIENT: AMEC Geomatix

PROJECT: FRP Shoreline Invest

SDG: TN43

CLIENT ID	ARI ID	DIL.	TIME	%R	AG	AL	AS	B	EA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	NO	NA	NI	PE	SB	SE	SI	SN	TI	TL	U	V	ZN
FRP-092111-001	TN43RA	1.00	13154																														
FRP-092111-001D	TN43RADUP	1.00	13162																														
FRP-092111-001S	TN43RASPX	1.00	13210																														
CCV	ACCV4	1.00	13235																														
CCB	CCB4	1.00	13263																														

General Chemistry Analysis
Report and Summary QC Forms

ARI Job ID: TN42, TN43

INORGANICS ANALYSIS DATA SHEET
pH by Method EPA 150.1

**ANALYTICAL
RESOURCES
INCORPORATED**

Data Release Authorized: *[Signature]*
 Reported: 09/23/11
 Date Received: 09/21/11
 Page 1 of 1

QC Report No: TN42-AMEC Geomatrix
 Project: FRP 2011 Shoreline Investigation
 8769

Client/ ARI ID	Date Sampled	Matrix	Analysis Date & Batch	RL	Result
FRP-092111-001 TN42A 11-20663	09/21/11	Water	09/21/11 17:28 092111#1	0.01	9.76
FRP-092111-002 TN42B 11-20664	09/21/11	Water	09/21/11 17:28 092111#1	0.01	7.93
FRP-092111-003 TN42C 11-20665	09/21/11	Water	09/21/11 17:28 092111#1	0.01	6.73
FRP-092111-004 TN42D 11-20666	09/21/11	Water	09/21/11 17:28 092111#1	0.01	6.70
FRP-092111-005 TN42E 11-20667	09/21/11	Water	09/21/11 17:28 092111#1	0.01	6.86
FRP-092111-006 TN42F 11-20668	09/21/11	Water	09/21/11 17:28 092111#1	0.01	6.91

Reported in std units

RL-Analytical reporting limit
 U-Undetected at reported detection limit

REPLICATE RESULTS-CONVENTIONALS
TN42-AMEC Geomatrix

ANALYTICAL
RESOURCES
INCORPORATED

Matrix: Water
Data Release Authorized: *[Signature]*
Reported: 09/23/11

Project: FRP 2011 Shoreline Investiga
Event: 8769
Date Sampled: 09/21/11
Date Received: 09/21/11

Analyte	Date	Units	Sample	Replicate(s)	RPD/RSD
ARI ID: TN42A Client ID: FRP-092111-001					
pH	09/21/11	std units	9.76	9.77	0.01

pH is evaluated as the Absolute Difference between the values rather than
Relative Percent Difference

LAB CONTROL RESULTS-CONVENTIONALS
TN42-AMEC Geomatrix

ANALYTICAL
RESOURCES
INCORPORATED

Matrix: Water
Data Release Authorized: ✓
Reported: 09/23/11

Project: FRP 2011 Shoreline Investiga
Event: 8769
Date Sampled: NA
Date Received: NA

Analyte	Date/Time	Units	LCS	Spike Added	Recovery
pH	09/21/11 std	units	6.97	7.00	0.03

pH is evaluated as the Absolute Difference between the values rather than Percent Recovery.